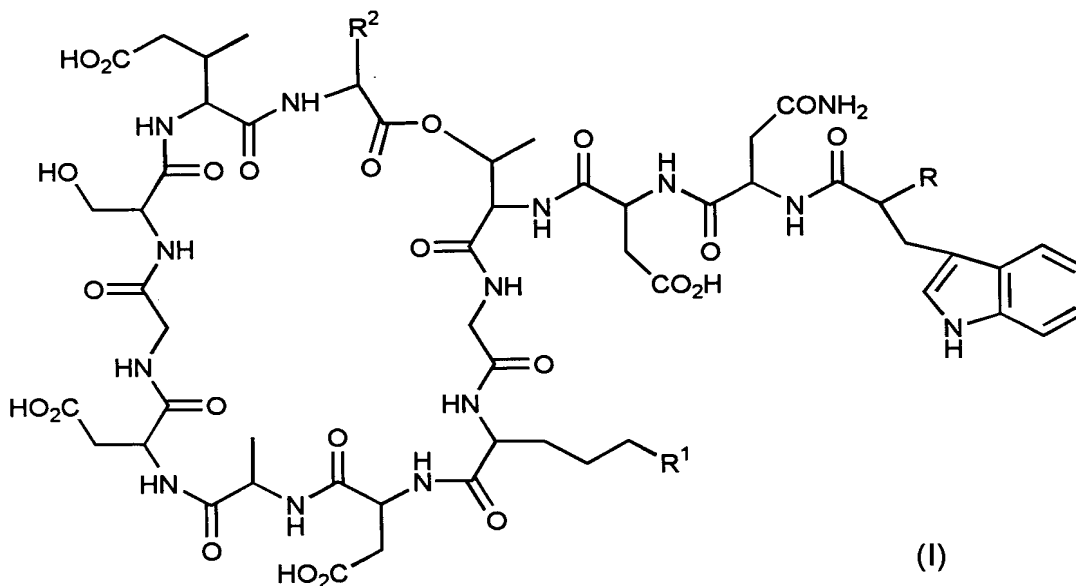
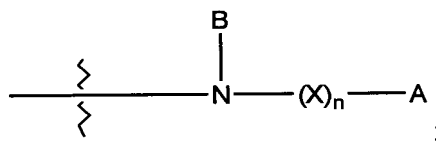


1. A compound having the formula (I):



wherein R is:



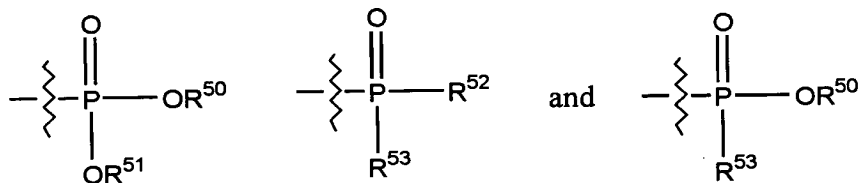
wherein n is 0 or 1;

wherein B is X^Y , H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein A is H, NH₂, NHR^A, NR^AR^B, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^A and R^B are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

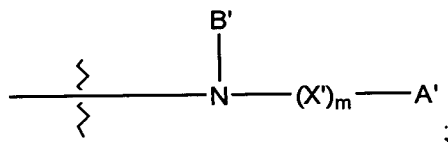
wherein when n is 0, then A is additionally selected from:



wherein each of R^{50} - R^{53} is independently selected from C_1 - C_{15} alkyl;

alternatively, wherein B and A together form a 5-7 membered heterocyclic or heteroaryl ring;

wherein R^1 is



wherein X' and X''' are independently selected from $C=O$, $C=S$, $C=NH$, $C=NR^{X'}$, $S=O$ or SO_2 ;

wherein m is 0 or 1;

wherein $R^{X'}$ is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

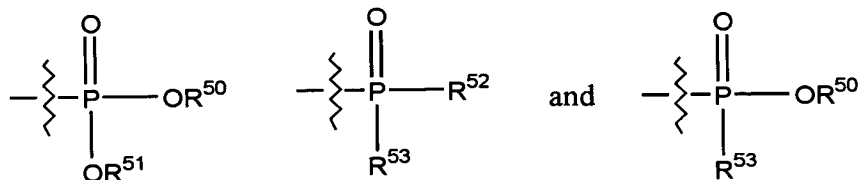
wherein B' is $X'''R^{Y'}$, H , alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein $R^{Y'}$ is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A' is H , NH_2 , $NHR^{A'}$, $NR^{A'}R^{B'}$, heteroaryl, cycloalkyl or heterocyclyl;

wherein $R^{A'}$ and $R^{B'}$ are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

wherein when m is 0, then A' is additionally selected from:



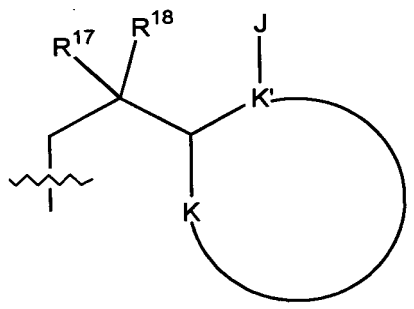
wherein each of R^{50} - R^{53} is independently selected from C_1 - C_{15} alkyl;
 provided that when B' is H and X' is $C=O$, then A' is other than
 (a) a pyridinyl ring substituted with one substituent $NHC(O)R^D$ or
 (b) a C_5 - C_6 saturated cycloalkyl ring substituted with one substituent

$NHC(O)R^D$;

wherein R^D is C_1 - C_{17} unsubstituted alkyl or C_2 - C_{17} unsubstituted alkenyl; and

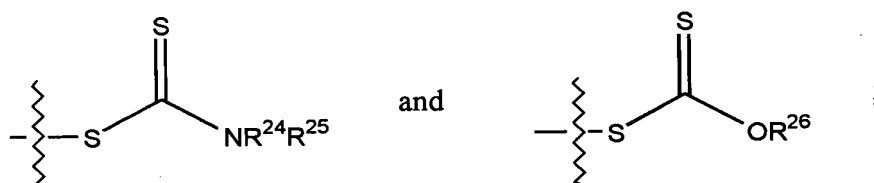
when B' is H and $m=0$, then A' is not H;

wherein R^2 is



wherein K and K' together form a C_3 - C_7 cycloalkyl or heterocyclyl ring or a C_5 - C_{10} aryl or heteroaryl ring;

wherein J is selected from the group consisting of hydrido, amino, NHR^J , $NR^J R^K$, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl, heterocyclyl, alkylamino, hydroxyl, thio, alkylthio, alkenylthio, sulfinyl, sulfonyl, azido, cyano, halo,



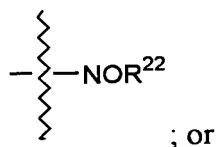
wherein each of R^{24} , R^{25} , and R^{26} is independently selected from the group consisting of alkyl, cycloalkyl, heterocyclyl, aryl and heteroaryl; or R^{24} and R^{25} together form a 5-8 membered heterocyclyl ring;

wherein R^J and R^K are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl; or

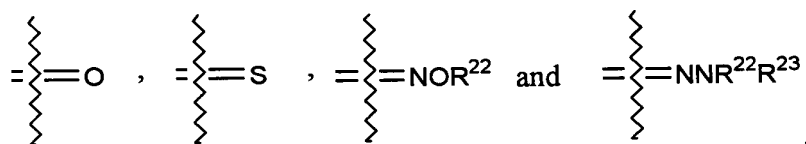
alternatively, wherein J, together with R^{17} , forms a 5-8 membered heterocyclyl or cycloalkyl ring; or

alternatively, wherein J, together with both R¹⁷ and R¹⁸, forms a 5-8 membered aryl, cycloalkyl, heterocyclyl or heteroaryl ring; and

wherein each of R¹⁷ and R¹⁸ is independently selected from the group consisting of hydrido, halo, hydroxyl, alkoxy, amino, thio, sulfinyl, sulfonyl and

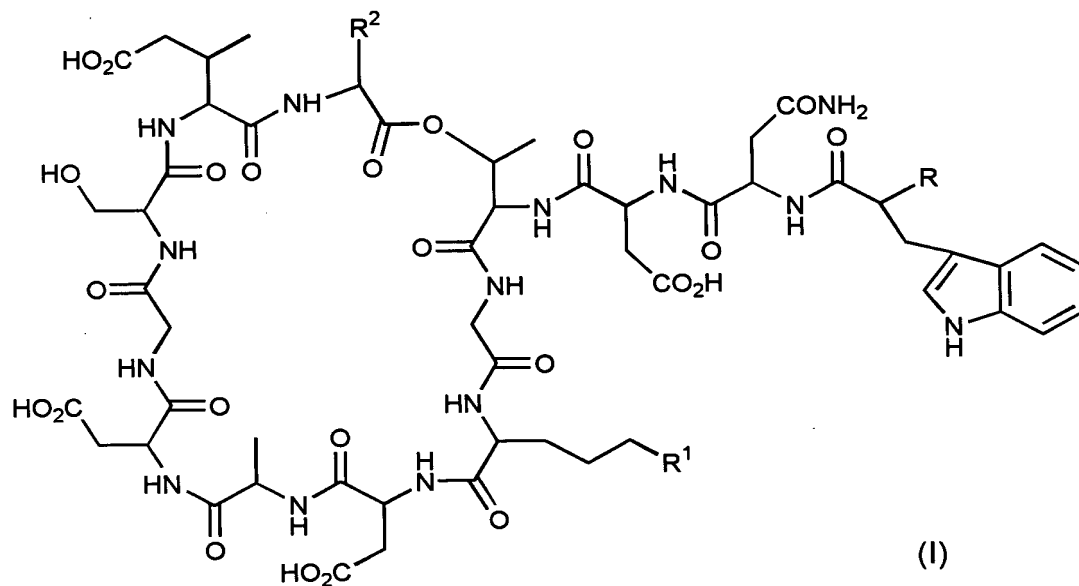


wherein R¹⁷ and R¹⁸ taken together can form a group consisting of ketal, thioketal,



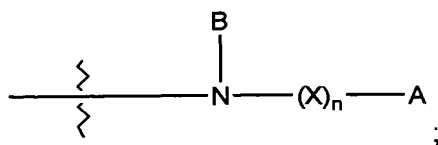
wherein each of R^{22} and R^{23} is independently selected from the group consisting of hydrido and alkyl.

2. A compound having the formula (I):



and salts thereof;

wherein R is:



wherein X and X" are independently selected from C=O, C=S, C=NH, C=NR^X, S=O or SO₂;

wherein n is 0 or 1;

wherein R^X is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

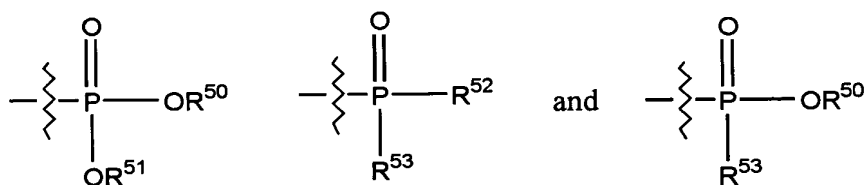
wherein B is X"R^Y, H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^Y is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A is H, NH₂, NHR^A, NR^AR^B, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^A and R^B are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

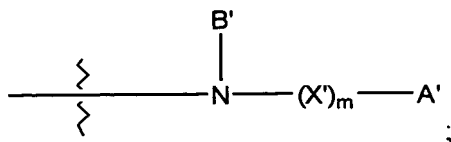
wherein when n is 0, then A is additionally selected from:



wherein each of R⁵⁰-R⁵³ is independently selected from C₁-C₁₅ alkyl;

alternatively, wherein B and A together form a 5-7 membered heterocyclic or heteroaryl ring;

wherein R¹ is



wherein X' and X''' are independently selected from C=O, C=S, C=NH, C=NR^X, S=O or SO₂;

wherein m is 0 or 1;

wherein $R^{X'}$ is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

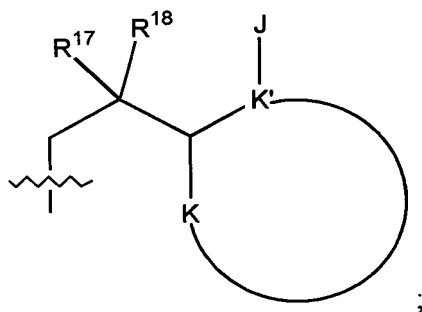
wherein B' is $X'''R^{Y'}$, H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein $R^{Y'}$ is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A' is aryl;

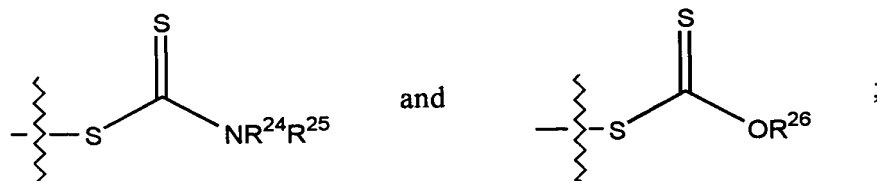
provided that when B' is H and X' is C=O, then A' is other than a phenyl ring substituted with substituent $NHC(O)R^D$, wherein R^D is C_1 - C_{17} unsubstituted alkyl or C_2 - C_{17} unsubstituted alkenyl, wherein said phenyl ring may be further optionally substituted with 1-2 substituents independently selected from amino, nitro, C_1 - C_3 alkyl, hydroxyl, C_1 - C_3 alkoxy, halo, mercapto, C_1 - C_3 alkylthio, carbamyl or C_1 - C_3 alkyl carbamyl;

wherein R^2 is



wherein K and K' together form a C_3 - C_7 cycloalkyl or heterocyclyl ring or a C_5 - C_{10} aryl or heteroaryl ring;

wherein J is selected from the group consisting of hydrido, amino, NHR^J , NR^JR^K , alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl, heterocyclyl, alkylamino, hydroxyl, thio, alkylthio, alkenylthio, sulfinyl, sulfonyl, azido, cyano, halo,



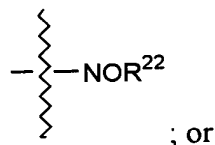
wherein each of R^{24} , R^{25} , and R^{26} is independently selected from the group consisting of alkyl, cycloalkyl, heterocyclyl, aryl and heteroaryl; or R^{24} and R^{25} together form a 5-8 membered heterocyclyl ring;

wherein R^J and R^K are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl; or

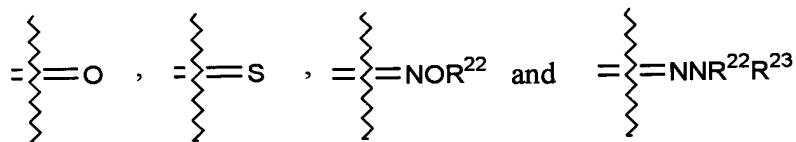
alternatively, wherein J, together with R^{17} , forms a 5-8 membered heterocyclyl or cycloalkyl ring; or

alternatively, wherein J, together with both R^{17} and R^{18} , forms a 5-8 membered aryl, cycloalkyl, heterocyclyl or heteroaryl ring; and

wherein each of R^{17} and R^{18} is independently selected from the group consisting of hydrido, halo, hydroxyl, alkoxy, amino, thio, sulfinyl, sulfonyl and

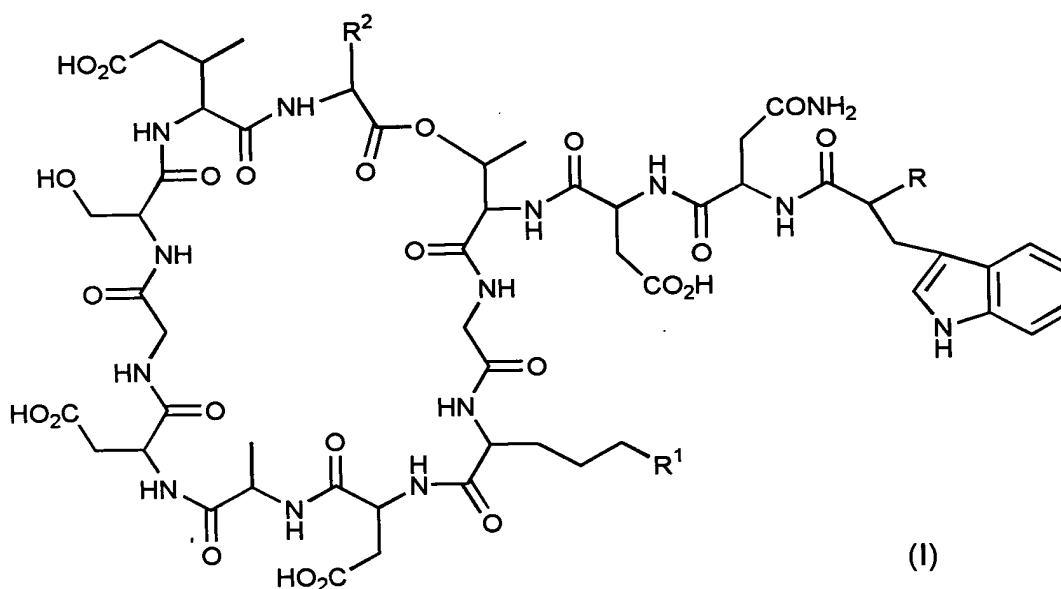


wherein R^{17} and R^{18} taken together can form a group consisting of ketal, thioketal,



wherein each of R^{22} and R^{23} is independently selected from the group consisting of hydrido and alkyl.

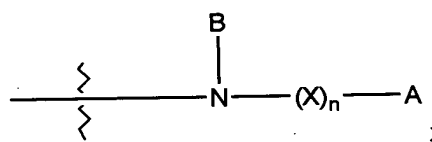
3. A compound having the formula (I):



(I)

and salts thereof;

wherein R is:



wherein X and X" are independently selected from C=O, C=S, C=NH, C=NR^X, S=O or SO₂;

wherein n is 0 or 1;

wherein R^X is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

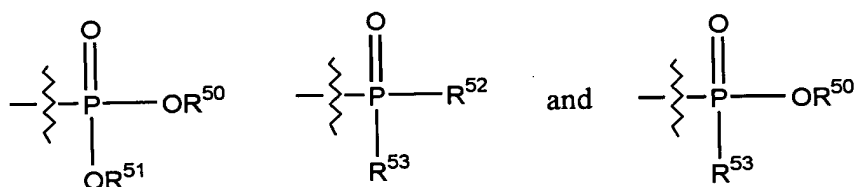
wherein B is X"R^Y, H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^Y is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A is H, NH₂, NHR^A, NR^AR^B, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^A and R^B are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

wherein when n is 0, then A is additionally selected from:

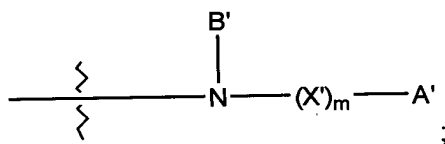


wherein each of R^{50} - R^{53} is independently selected from C_1 - C_{15} alkyl;

alternatively, wherein B and A together form a 5-7 membered

heterocyclic or heteroaryl ring;

wherein R^1 is



wherein X' and X'' are independently selected from $\text{C}=\text{O}$, $\text{C}=\text{S}$, $\text{C}=\text{NH}$, $\text{C}=\text{NR}^{\text{X}'}$, $\text{S}=\text{O}$ or SO_2 ;

wherein m is 0 or 1;

wherein $\text{R}^{\text{X}'}$ is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

wherein B' is $\text{X}''\text{R}^{\text{Y}'}$, H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein $\text{R}^{\text{Y}'}$ is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A' is alkyl, alkenyl, alkynyl, alkoxy or aryloxy;

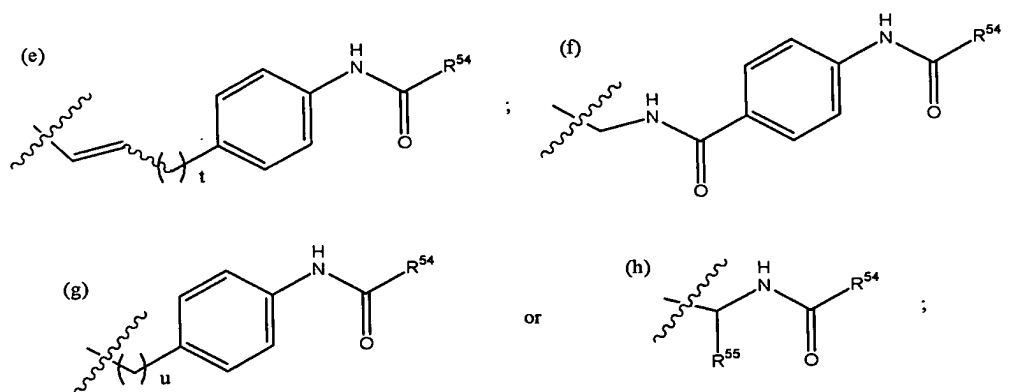
provided that when B' is H and X' is $\text{C}=\text{O}$, then A' is other than

(a) $-(\text{C}_1\text{-C}_{16}\text{ unsubstituted alkyl})\text{-NH}_2$;

(b) $-(\text{C}_1\text{-C}_{10}\text{ unsubstituted alkyl})\text{-NHC(O)R}^{\text{D}}$, wherein R^{D} is $-\text{C}_1\text{-C}_{18}$ alkyl, optionally substituted with up to one hydroxyl, carboxyl or $\text{C}_1\text{-C}_3$ alkoxy, or one to three halo substituents;

(c) $-\text{C}_1\text{-C}_{18}$ alkyl, optionally substituted with up to one hydroxyl, carboxyl or $\text{C}_1\text{-C}_3$ alkoxy, or one to three halo substituents;

(d) $-\text{C}_4\text{-C}_{18}$ unsubstituted alkenyl;

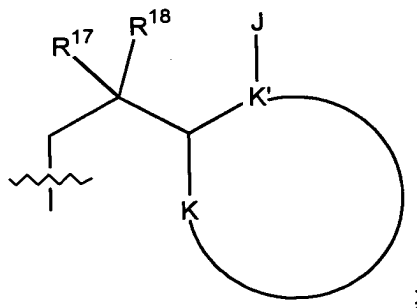


wherein R^{54} is selected from $\text{C}_1\text{-C}_{17}$ -unsubstituted alkyl or $\text{C}_2\text{-C}_{17}$ -unsubstituted alkenyl; wherein R^{55} is selected from hydroxyethyl, hydroxymethyl, mercaptomethyl, mercaptoethyl, methylthioethyl, 2-thienyl, 3-indolemthyl, phenyl optionally substituted with a group selected from halo, nitro, $\text{C}_1\text{-C}_3$ -unsubstituted alkyl, hydroxy, $\text{C}_1\text{-C}_3$ -unsubstituted alkoxy, $\text{C}_1\text{-C}_3$ -unsubstituted alkylthio, carbamyl or $\text{C}_1\text{-C}_3$ unsubstituted alkylcarbamyl; or benzyl optionally substituted with a group selected from halo, nitro, $\text{C}_1\text{-C}_3$ -unsubstituted alkyl, hydroxy, $\text{C}_1\text{-C}_3$ -unsubstituted alkoxy, $\text{C}_1\text{-C}_3$ -unsubstituted alkylthio, carbamyl or $\text{C}_1\text{-C}_3$ unsubstituted alkylcarbamyl; wherein t is 0 or 1 and wherein u is an integer from 1-3; and

when B is H and X is $\text{C}=\text{O}$, then X , together with A , does not form a carbamate amino protecting group; and

wherein when B' is H and m is 0, then A' is other than $\text{C}_4\text{-C}_{14}$ unsubstituted alkyl;

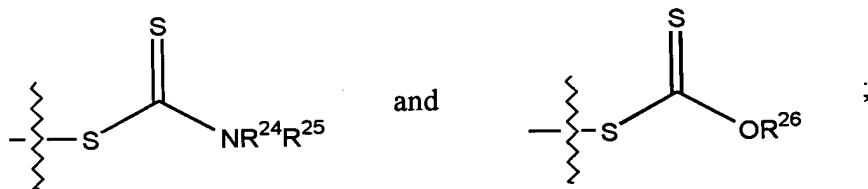
wherein R^2 is



wherein K and K' together form a $\text{C}_3\text{-C}_7$ cycloalkyl or heterocyclyl ring or a $\text{C}_5\text{-C}_{10}$ aryl or heteroaryl ring;

wherein J is selected from the group consisting of hydrido, amino, NHR^J , NR^JR^K , alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl,

heterocyclyl, alkylamino, hydroxyl, thio, alkylthio, alkenylthio, sulfinyl, sulfonyl, azido, cyano, halo,



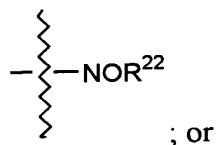
wherein each of R^{24} , R^{25} , and R^{26} is independently selected from the group consisting of alkyl, cycloalkyl, heterocyclyl, aryl and heteroaryl; or R^{24} and R^{25} together form a 5-8 membered heterocyclyl ring;

wherein R^J and R^K are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl; or

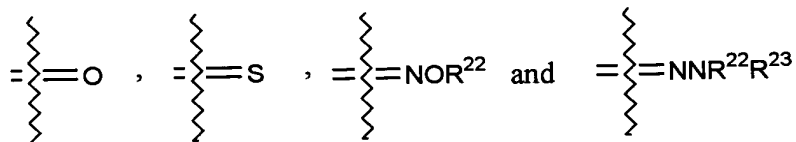
alternatively, wherein J, together with R^{17} , forms a 5-8 membered heterocyclyl or cycloalkyl ring; or

alternatively, wherein J, together with both R^{17} and R^{18} , forms a 5-8 membered aryl, cycloalkyl, heterocyclyl or heteroaryl ring; and

wherein each of R^{17} and R^{18} is independently selected from the group consisting of hydrido, halo, hydroxyl, alkoxy, amino, thio, sulfinyl, sulfonyl and

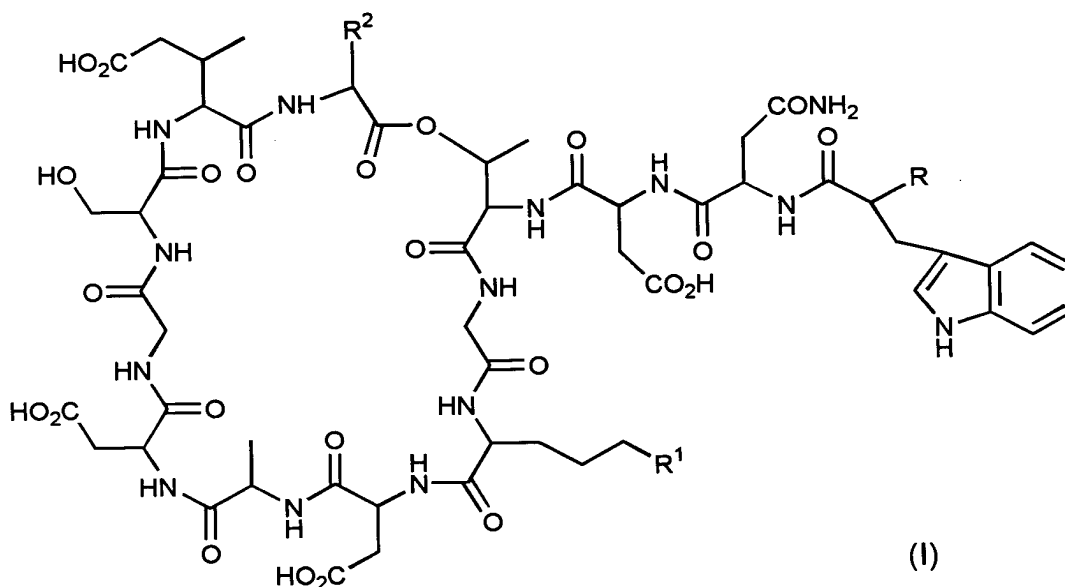


wherein R^{17} and R^{18} taken together can form a group consisting of ketal, thioketal,



wherein each of R^{22} and R^{23} is independently selected from the group consisting of hydrido and alkyl.

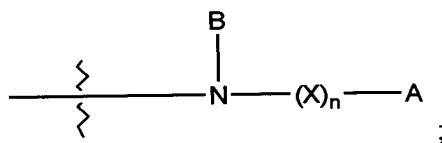
4. A compound having the formula (I):



(I)

and salts thereof;

wherein R is:



wherein X and X" are independently selected from C=O, C=S, C=NH, C=NR^X, S=O or SO₂;

wherein n is 0 or 1;

wherein R^X is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

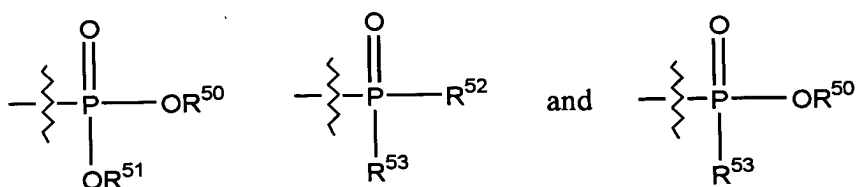
wherein B is X"R^Y, H, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^Y is selected from hydrido, alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or hydroxyl;

wherein A is H, NH₂, NHR^A, NR^AR^B, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl or heterocyclyl;

wherein R^A and R^B are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

wherein when n is 0, then A is additionally selected from:

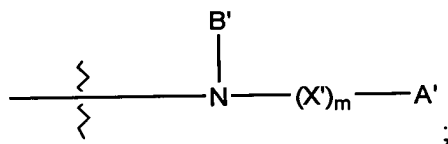


wherein each of R^{50} - R^{53} is independently selected from C_1 - C_{15} alkyl;

alternatively, wherein B and A together form a 5-7 membered

heterocyclic or heteroaryl ring;

wherein R^1 is



wherein X' and X'' are independently selected from $\text{C}=\text{O}$, $\text{C}=\text{S}$, $\text{C}=\text{NH}$, $\text{C}=\text{NR}^{\text{X}'}$, $\text{S}=\text{O}$ or SO_2 ;

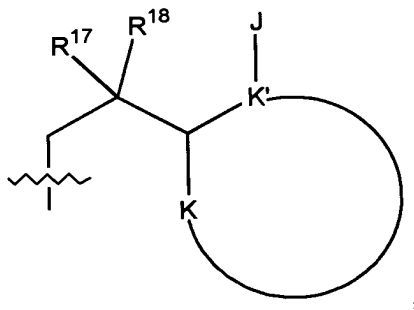
wherein m is 0 or 1;

wherein $\text{R}^{\text{X}'}$ is selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl, hydroxyl, alkoxy, carboxy or carboalkoxy;

wherein B' and A' together form a 5-7 membered heterocyclic or heteroaryl ring;

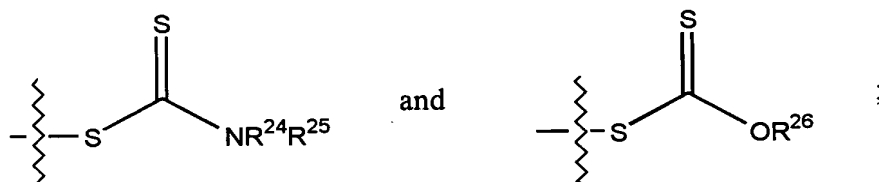
wherein $\text{R}^{\text{A}'}$ and $\text{R}^{\text{B}'}$ are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl, heterocyclyl or carboalkoxy;

wherein R^2 is



wherein K and K' together form a C_3 - C_7 cycloalkyl or heterocyclyl ring or a C_5 - C_{10} aryl or heteroaryl ring;

wherein J is selected from the group consisting of hydrido, amino, NHR^{J} , $\text{NR}^{\text{J}}\text{R}^{\text{K}}$, alkyl, alkenyl, alkynyl, alkoxy, aryloxy, aryl, heteroaryl, cycloalkyl, heterocyclyl, alkylamino, hydroxyl, thio, alkylthio, alkenylthio, sulfinyl, sulfonyl, azido, cyano, halo,



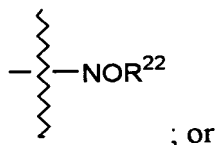
wherein each of R²⁴, R²⁵, and R²⁶ is independently selected from the group consisting of alkyl, cycloalkyl, heterocyclyl, aryl and heteroaryl; or R²⁴ and R²⁵ together form a 5-8 membered heterocyclyl ring;

wherein R^J and R^K are independently selected from alkyl, alkenyl, alkynyl, aryl, heteroaryl, cycloalkyl or heterocyclyl; or

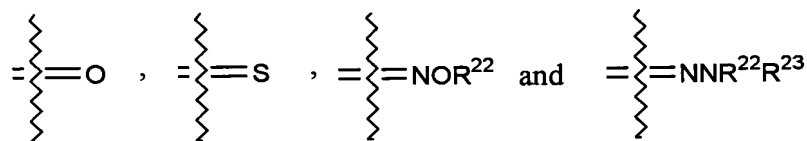
alternatively, wherein J, together with R¹⁷, forms a 5-8 membered heterocyclyl or cycloalkyl ring; or

alternatively, wherein J, together with both R¹⁷ and R¹⁸, forms a 5-8 membered aryl, cycloalkyl, heterocyclyl or heteroaryl ring; and

wherein each of R¹⁷ and R¹⁸ is independently selected from the group consisting of hydrido, halo, hydroxyl, alkoxy, amino, thio, sulfinyl, sulfonyl and

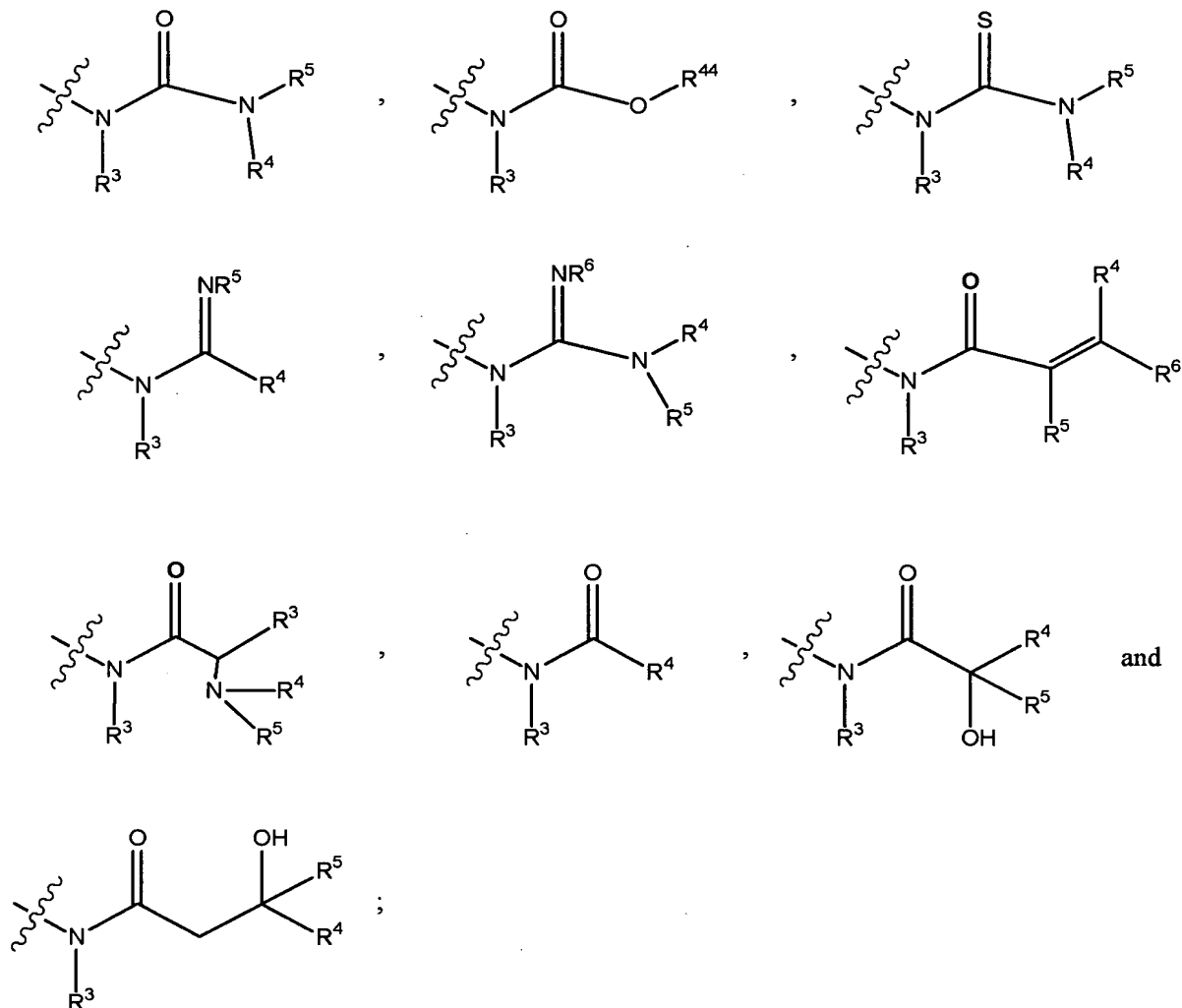


wherein R¹⁷ and R¹⁸ taken together can form a group consisting of ketal, thioketal,



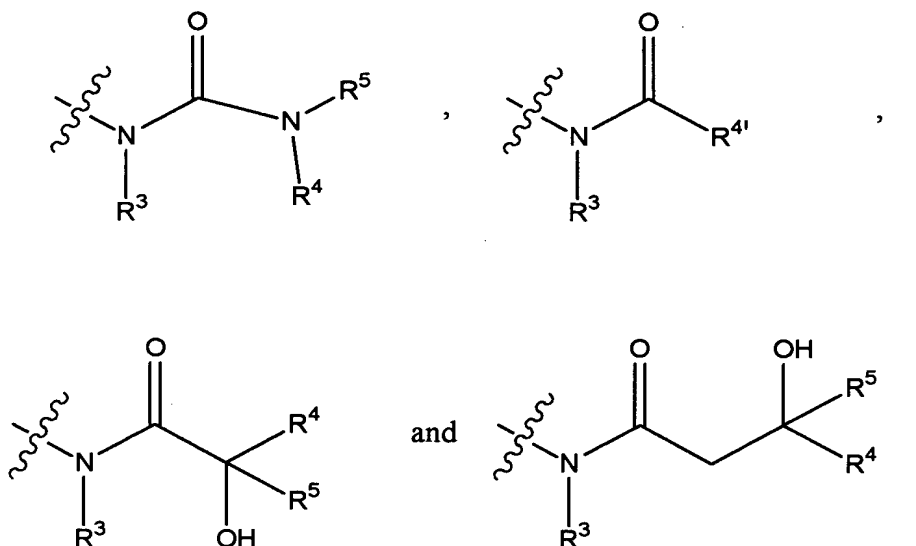
wherein each of R²² and R²³ is independently selected from the group consisting of hydrido and alkyl.

5. The compound according to any of claims 1-4, wherein R is selected from the group consisting of:



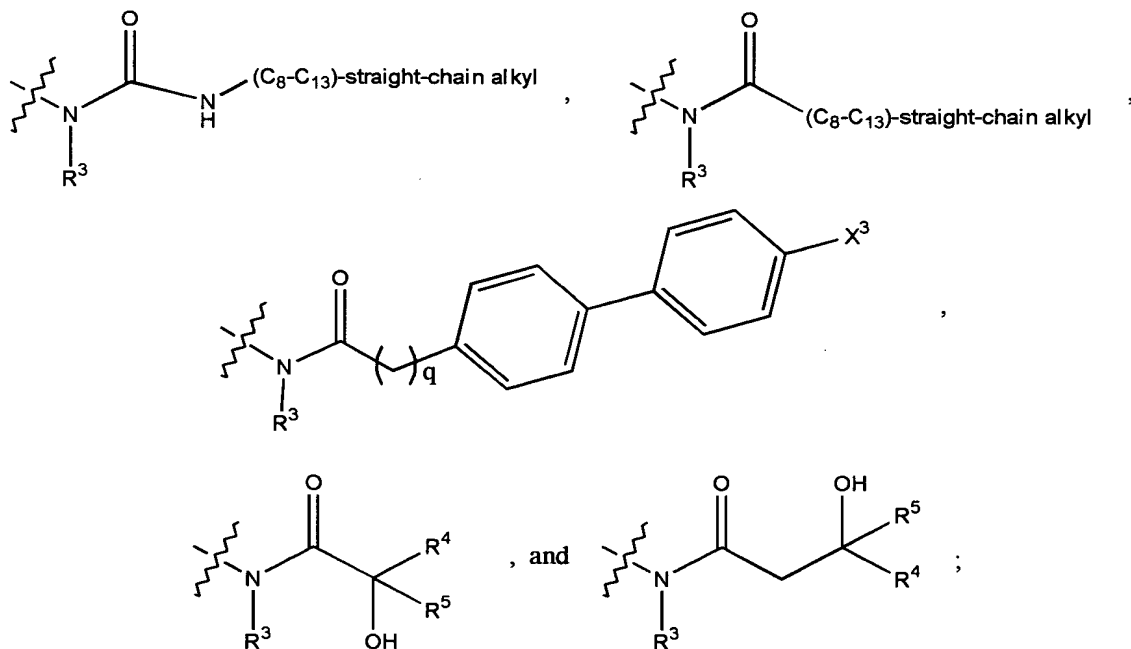
wherein each of R^3 , R^4 , R^5 , and R^6 is independently selected from the group consisting of hydrido, alkyl, aryl, heterocyclyl and heteroaryl, and wherein R^{44} is selected from the group consisting of alkyl, aryl, heterocyclyl and heteroaryl.

6. The compound according to claim 5, wherein R is selected from



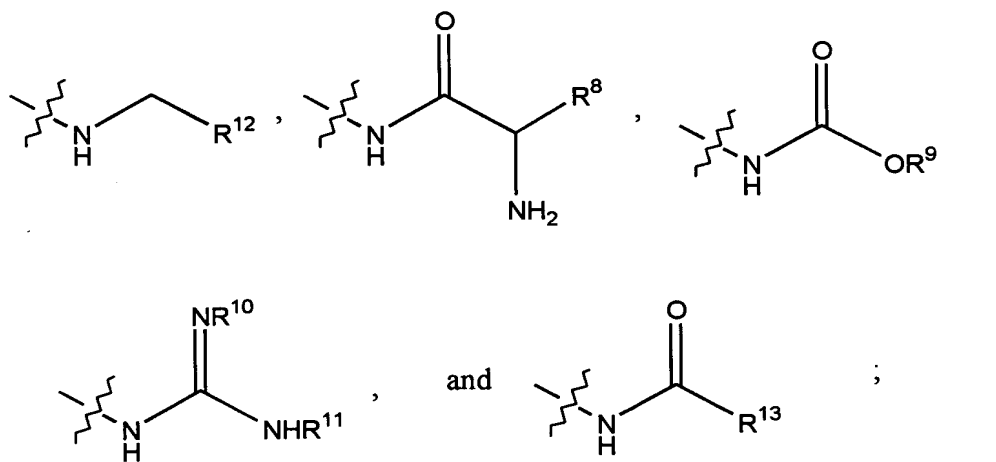
wherein R^{41} is selected from the group consisting of alkyl, aryl-substituted alkyl, substituted phenyl, heteroaryl, heterocyclyl, optionally substituted (C_8-C_{14}) -straight chain alkyl and ---SR^7 ; wherein R^7 is an alkyl group.

7. The compound according to claim 6, wherein R is selected from the group consisting of



wherein X^3 is chloro or trifluoromethyl and wherein q is 0 or 1.

8. The compound according to any of claims 1- 4, wherein R^1 is selected from the group consisting of:



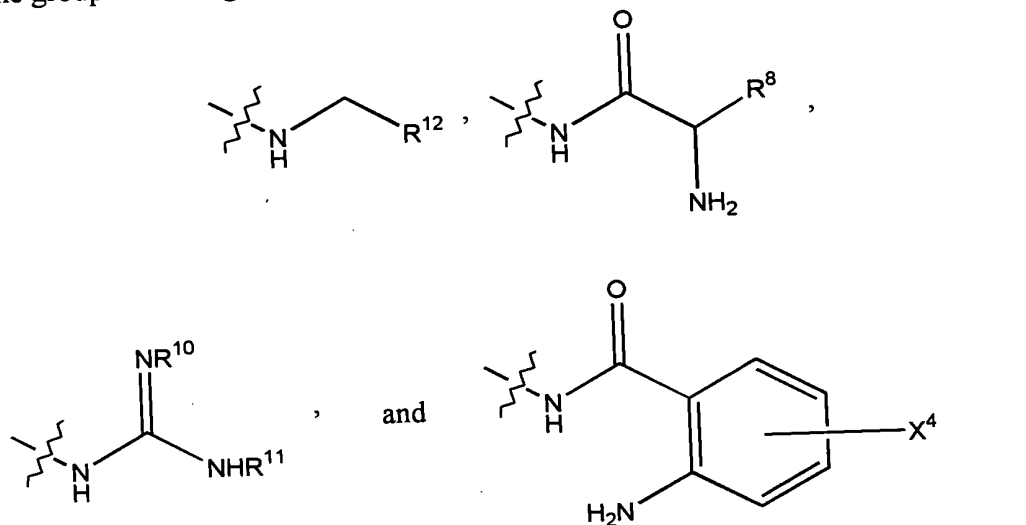
wherein R^8 is selected from a natural amino acid side chain or an amino acid side chain that is not naturally occurring;

wherein each of R^9 , R^{10} and R^{11} is selected from hydrido, alkyl, aryl, heterocyclyl and heteroaryl;

wherein R^{12} is selected from the group consisting of heterocyclyl, heteroaryl, aryl, and alkyl and

wherein R^{13} is selected from (C_1 - C_3 -alkyl) and aryl.

9. The compound according to claim 8, wherein R^1 is selected from the group consisting of:



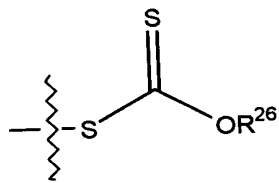
wherein R^8 is selected from tryptophan side chain and lysine side chain;

wherein each of R^{10} and R^{11} is independently selected from hydrido and alkyl;

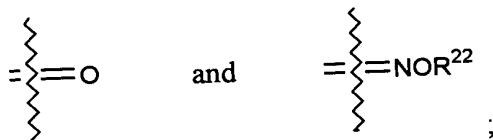
wherein R^{12} is selected from imidazolyl, N-methylimidazolyl, indolyl, quinoliny, benzyloxybenzyl, and benzylpiperidenylbenzyl; and

wherein X is selected from fluoro, and trifluoromethyl.

10. The compound according to any of claims 1-4, wherein J is selected from the group consisting of hydrido, amino, azido and



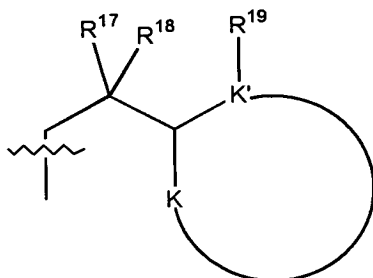
wherein R^{17} and R^{18} taken together form a group selected from ketal,



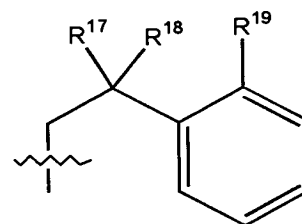
or wherein R^{17} is hydroxyl when R^{18} is hydrido;

or wherein J, together with R¹⁷, forms a heterocyclyl ring.

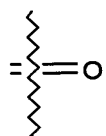
11. The compound according to claim 10, wherein R² is selected from the group consisting of



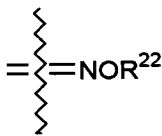
and



wherein R¹⁷ and R¹⁸ taken together form a group selected from

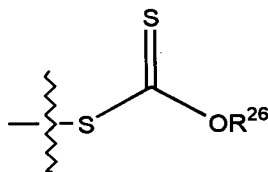


and



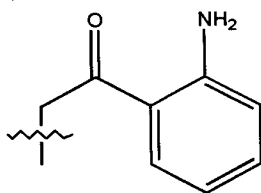
, wherein R²² is selected from the group

consisting of H and alkyl; and wherein R¹⁹ is selected from the group consisting of

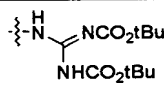
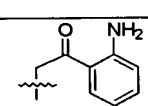
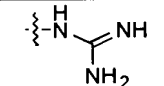
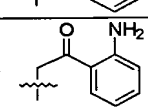
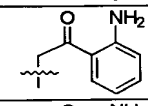
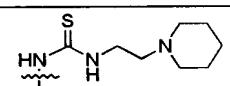
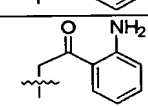
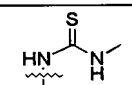
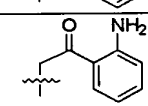
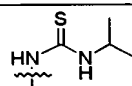
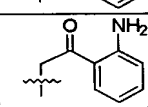
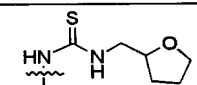
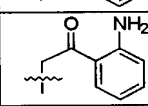
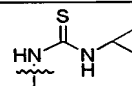
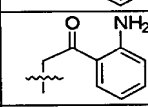
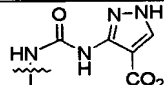
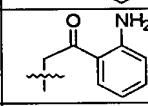
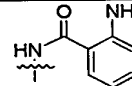
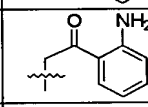
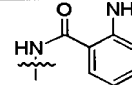
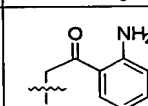
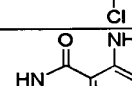
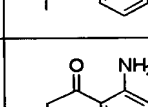
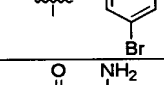
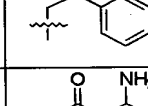
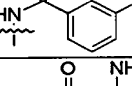
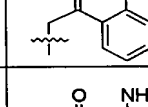
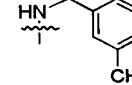
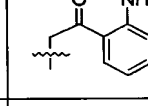
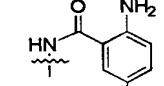
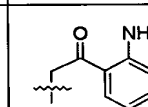


hydrido, amino, azido and

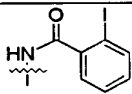
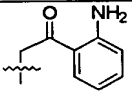
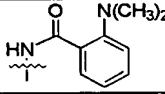
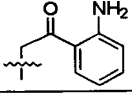
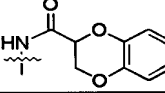
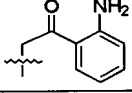
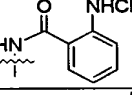
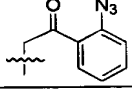
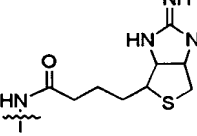
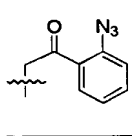
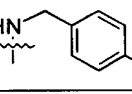
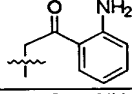
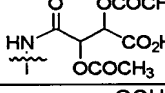
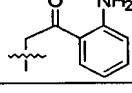
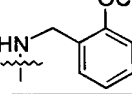
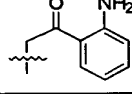
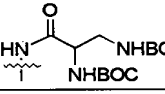
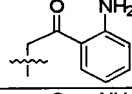
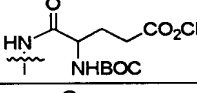
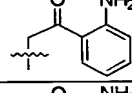
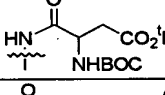
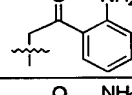
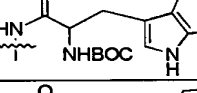
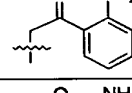
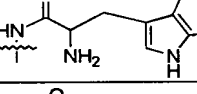
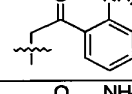
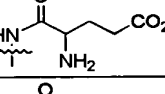
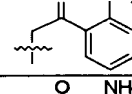
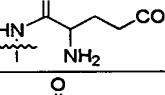
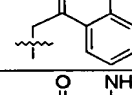
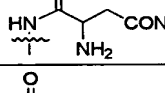
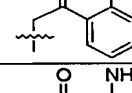
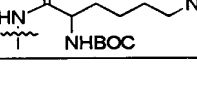
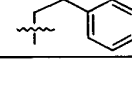
12. The compound according to claim 11, wherein R² is



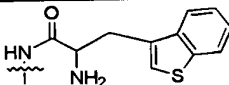
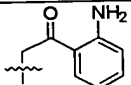
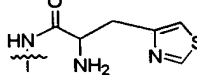
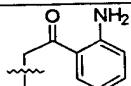
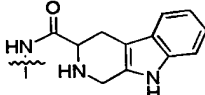
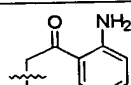
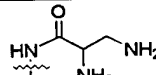
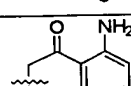
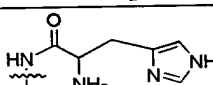
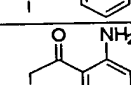
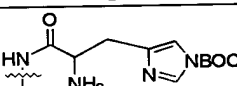
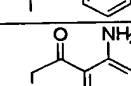
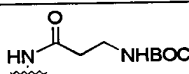
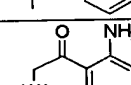
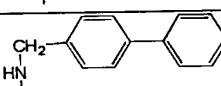
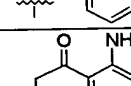
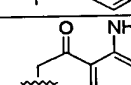
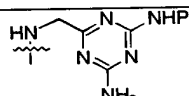
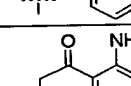
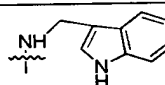
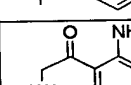
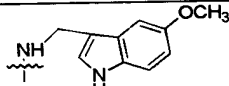
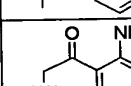
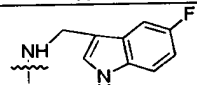
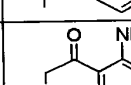
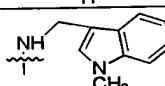
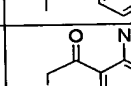
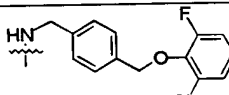
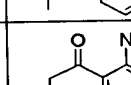
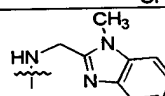
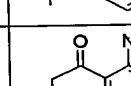

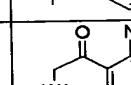
13. The compound according to any one of claims 1-4 wherein said compound is selected from

Cpd #	R	R ¹	R ²
1	NHCO(CH ₂) ₈ CH ₃		
2	NHCO(CH ₂) ₈ CH ₃		
3	NHCO(CH ₂) ₈ CH ₃	NHSO ₂ Ph	
4	NHCO(CH ₂) ₈ CH ₃		
5	NHCO(CH ₂) ₈ CH ₃		
6	NHCO(CH ₂) ₈ CH ₃		
7	NHCO(CH ₂) ₈ CH ₃		
8	NHCO(CH ₂) ₈ CH ₃		
9	NHCO(CH ₂) ₈ CH ₃		
10	NHCO(CH ₂) ₈ CH ₃		
11	NHCO(CH ₂) ₈ CH ₃		
12	NHCO(CH ₂) ₈ CH ₃		
13	NHCO(CH ₂) ₈ CH ₃		
14	NHCO(CH ₂) ₈ CH ₃		
15	NHCO(CH ₂) ₈ CH ₃		
16	NHCO(CH ₂) ₈ CH ₃		

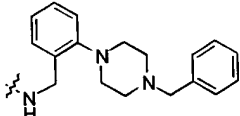
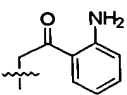
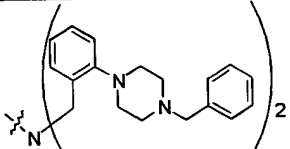
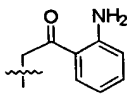
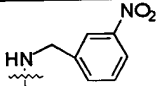
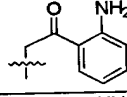
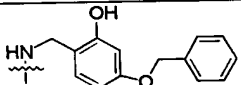
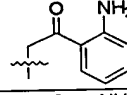
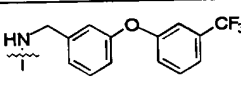
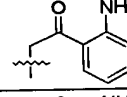
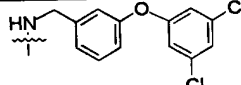
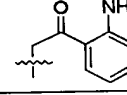
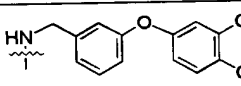
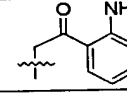
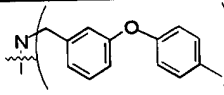
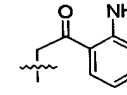
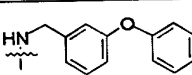
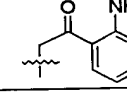
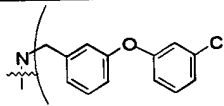
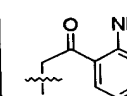
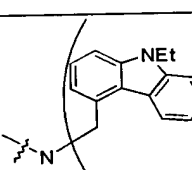
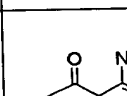
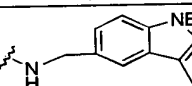
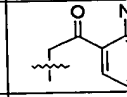

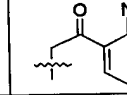
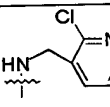
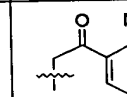
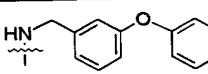
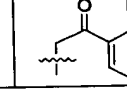
17	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
18	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
19	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
20	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
21	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
22	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
23	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
24	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
25	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
26	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
27	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
28	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
29	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
30	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
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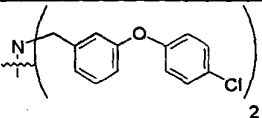
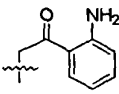
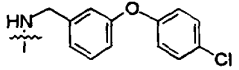
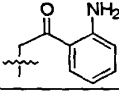
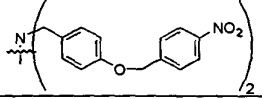
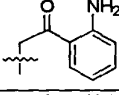
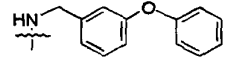
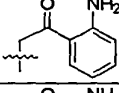
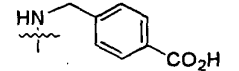
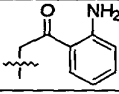
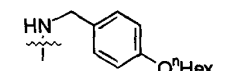
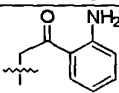
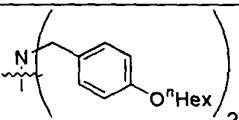
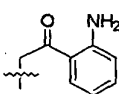
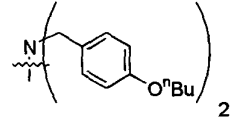
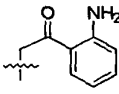
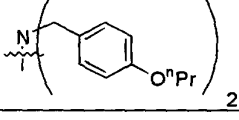
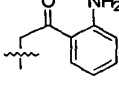
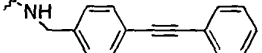
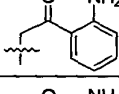
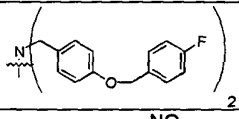
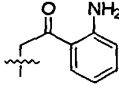
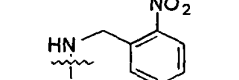
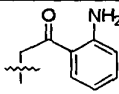
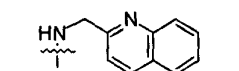
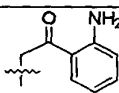
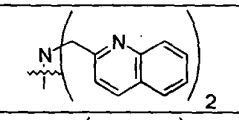
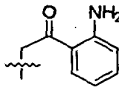
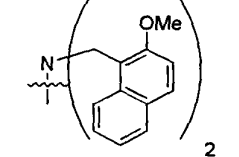
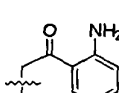
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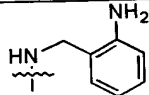
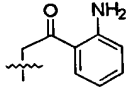
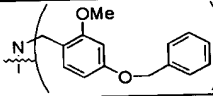
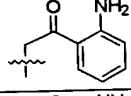
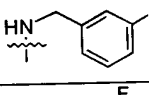
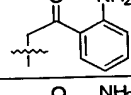
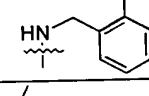
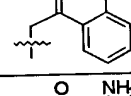
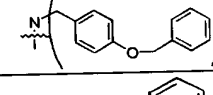
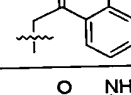
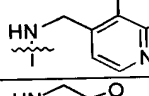
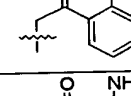
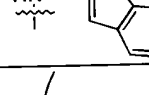
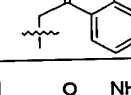
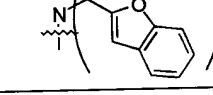
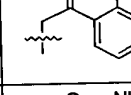
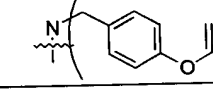
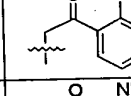
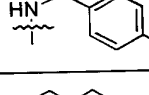
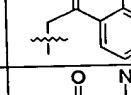
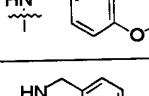
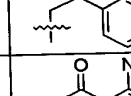
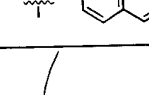
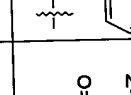
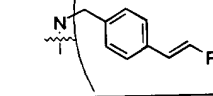
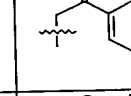
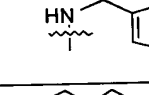
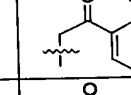
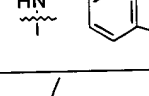
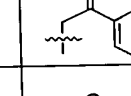
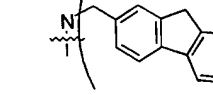
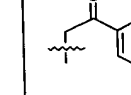
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86			
87	NHCO(CH ₂) ₈ CH ₃		
88	NHCO(CH ₂) ₈ CH ₃		
89	NHCO(CH ₂) ₈ CH ₃		
90	NHCO(CH ₂) ₈ CH ₃		
91	NHCO(CH ₂) ₈ CH ₃		
92	NHCO(CH ₂) ₈ CH ₃		
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94	NHCO(CH ₂) ₈ CH ₃		
95	NHCO(CH ₂) ₈ CH ₃		
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98	NHCO(CH ₂) ₈ CH ₃		
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101			
102	NHCO(CH ₂) ₁₁ CH ₃		

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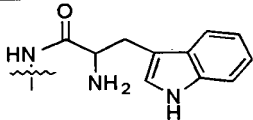
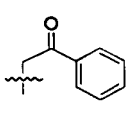
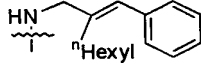
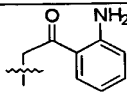
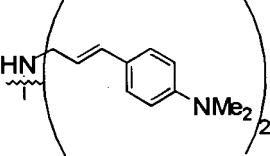
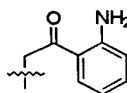
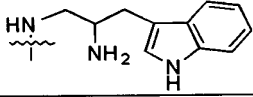
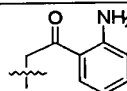
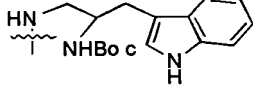
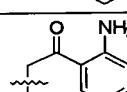
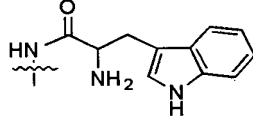
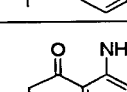
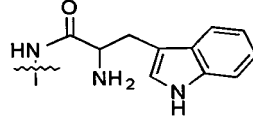
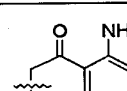
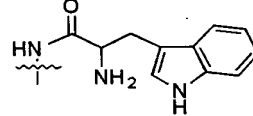
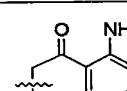
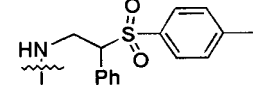
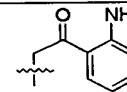
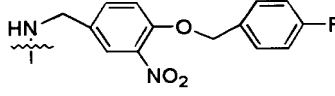
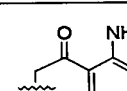
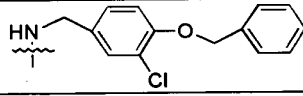
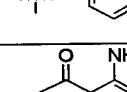
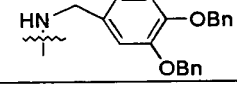
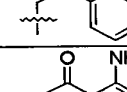
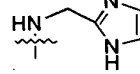
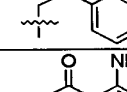
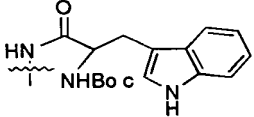
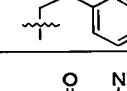
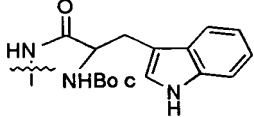
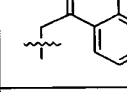
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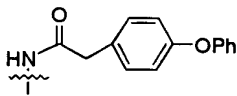
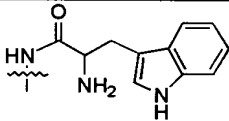
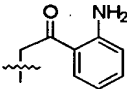
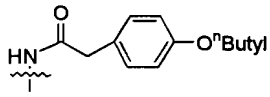
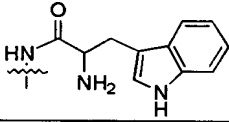
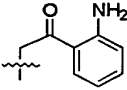
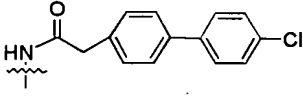
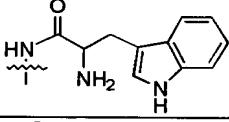
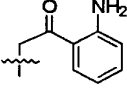
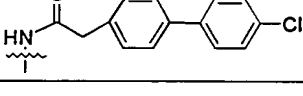
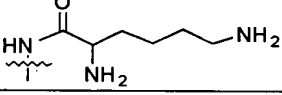
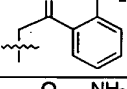
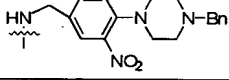
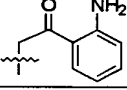
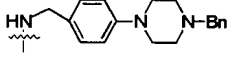
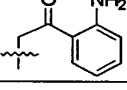
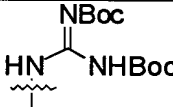
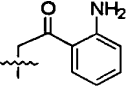
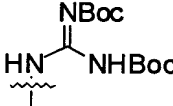
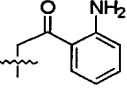
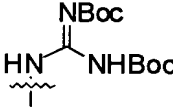
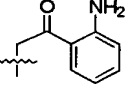
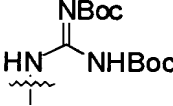
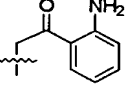
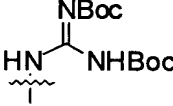
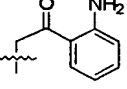
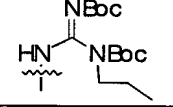
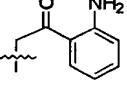
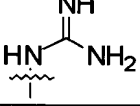
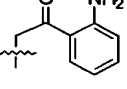
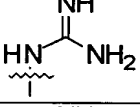
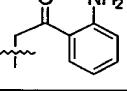
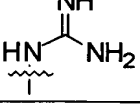
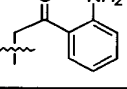
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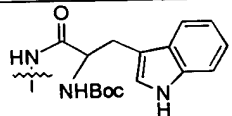
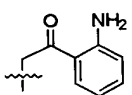
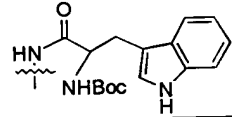
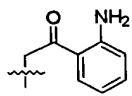
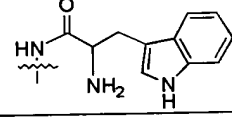
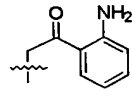
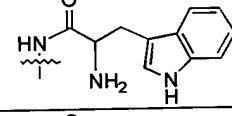
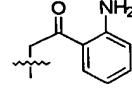
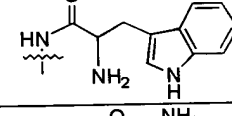
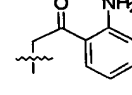
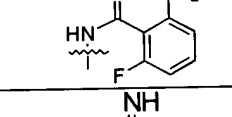
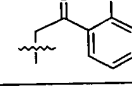
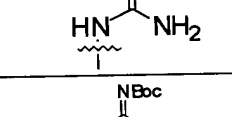
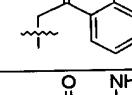
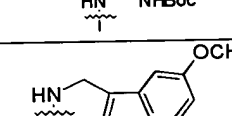
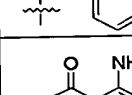
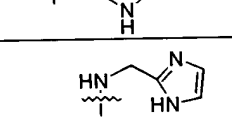
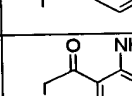
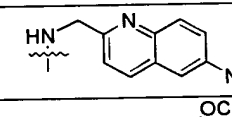
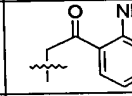
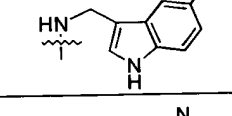
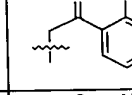
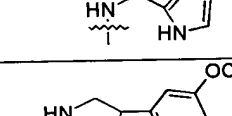
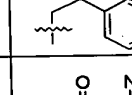
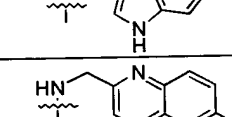
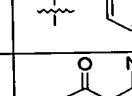
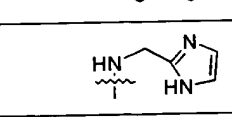
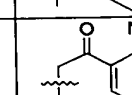


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233	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
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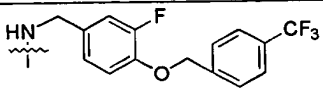
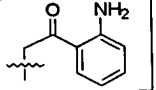
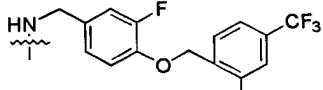
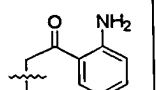
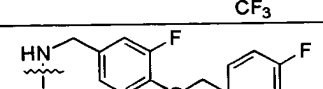
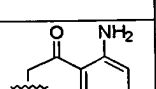
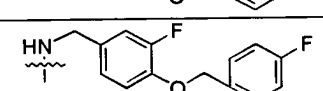
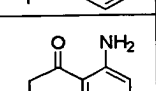
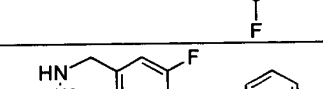
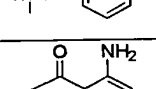
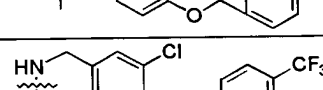
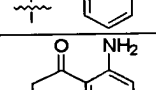
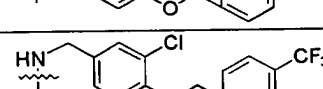
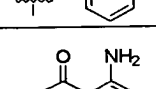
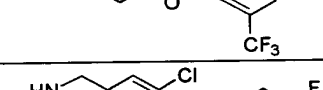
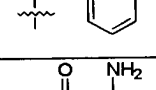
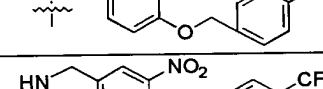
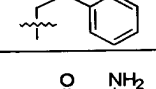
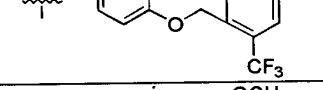
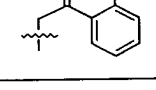
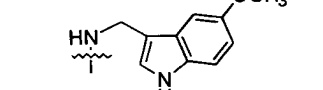
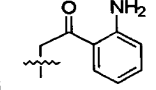
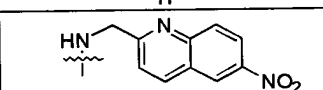
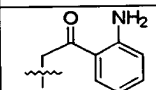
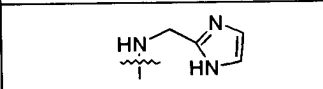
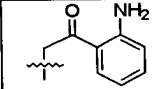
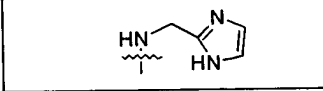
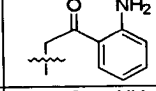
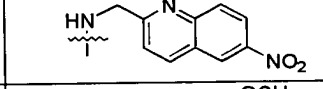
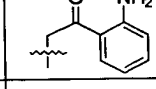
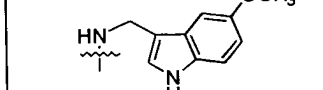
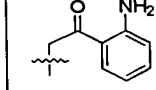
248			
249			
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262	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		

263			
264			
265	NHCO(CH ₂) ₈ CH ₃		
266	NHCO(CH ₂) ₈ CH ₃		
267	NHCO(CH ₂) ₈ CH ₃		
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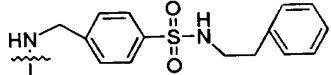
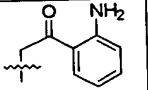
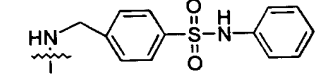
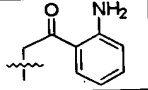
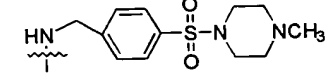
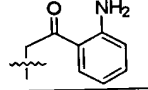
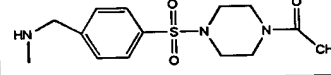
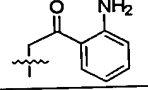
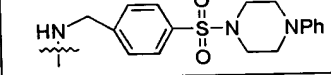
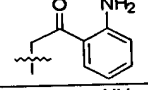
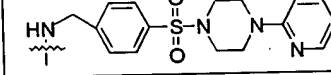
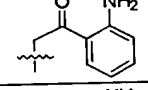
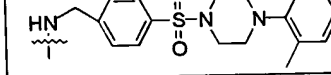
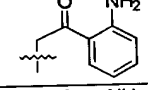
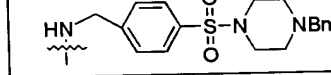
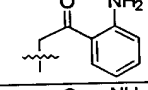
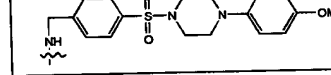
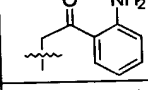
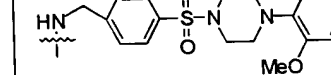
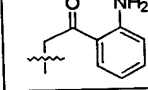
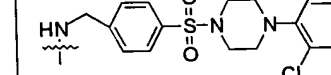
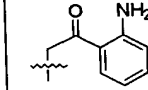
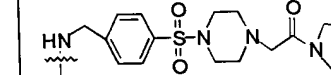
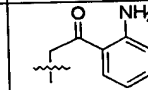
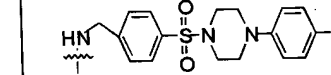
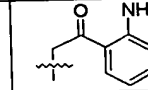
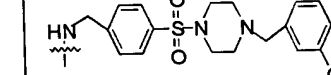
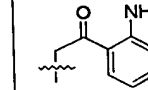
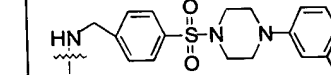
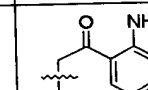
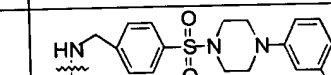
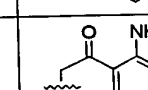
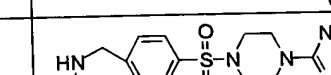
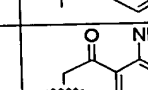
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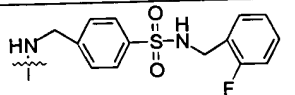
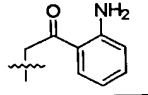
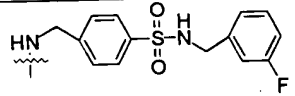
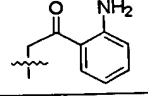
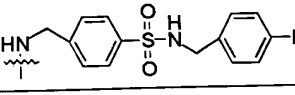
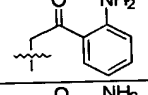
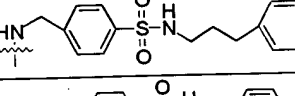
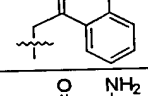
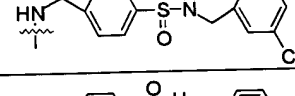
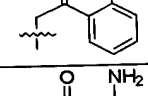
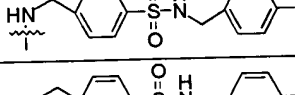
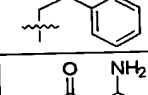
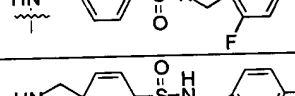
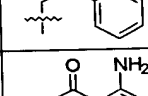
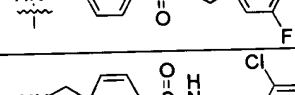
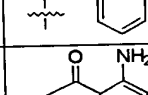
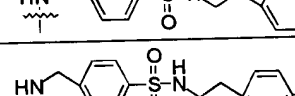
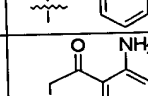
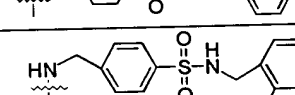
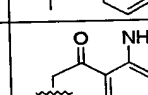
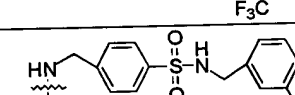
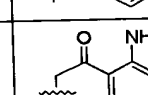
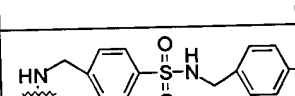
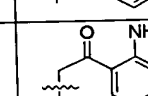
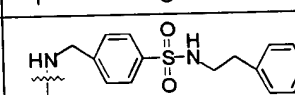
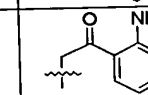
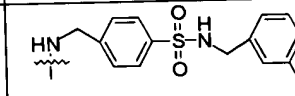
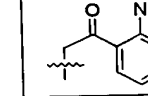
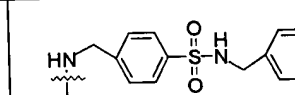
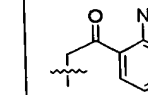
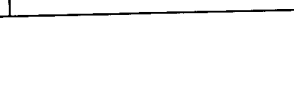

292			
293	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
294	$\text{NHCO}(\text{CH}_2)_7\text{CH}_3$		
295	$\text{NHCO}(\text{CH}_2)_{11}\text{CH}_3$		
296	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
297	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
298	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
299	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		
300	$\text{NHCONH}(\text{CH}_2)_{11}\text{CH}_3$		
301	$\text{NHCO}(\text{CH}_2)_{11}\text{CH}_3$		
302	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
303	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
304	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
305	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		
306	$\text{NHCONH}(\text{CH}_2)_{11}\text{CH}_3$		

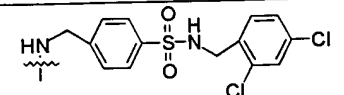
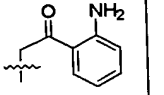
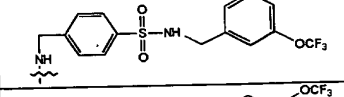
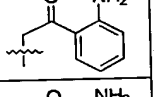
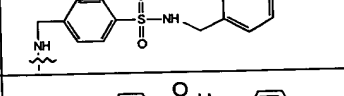
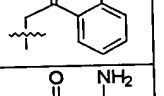
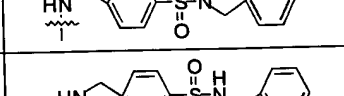
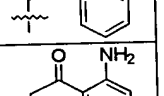
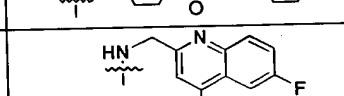
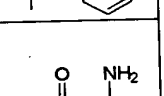
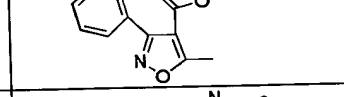
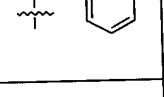
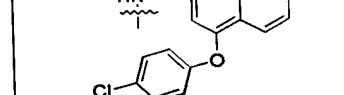
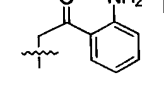
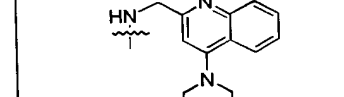
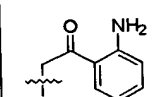
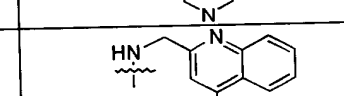
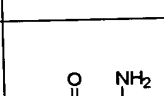
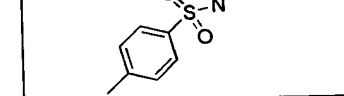
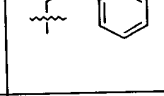
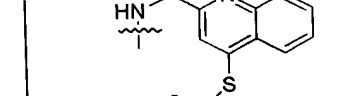
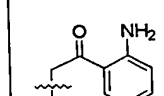
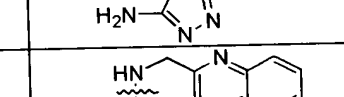
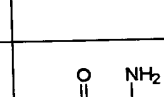
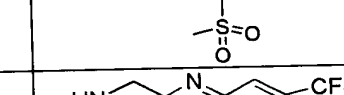
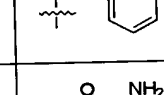
307	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
308	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
309	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
310	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
311	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
312	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
313	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
314	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		
315	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
316	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
317	$\text{NHCONH}(\text{CH}_2)_7\text{CH}_3$		
318	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
319	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
320	$\text{NHCO}(\text{CH}_2)_{11}\text{CH}_3$		
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322	$\text{NHCO}(\text{CH}_2)_{11}\text{CH}_3$		

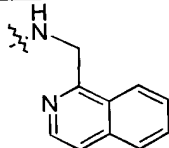
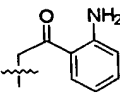
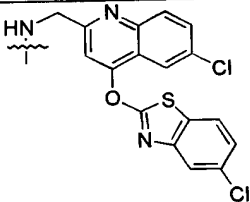
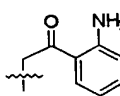
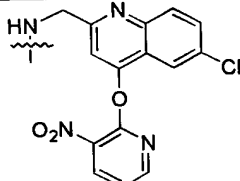
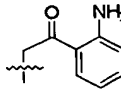
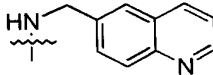
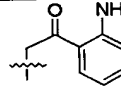
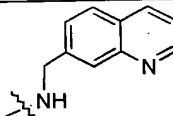
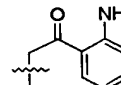
323	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
324	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
325	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
326	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
327	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
328	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
329	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
330	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
331	$\text{NHCO}(\text{CH}_2)_8\text{CH}_3$		
332	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
333	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
334	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
335	$\text{NHCONH}(\text{CH}_2)_{11}\text{CH}_3$		
336	$\text{NHCONH}(\text{CH}_2)_{11}\text{CH}_3$		
337	$\text{NHCONH}(\text{CH}_2)_{11}\text{CH}_3$		
338	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		

339	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
340	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
341	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
342	$\text{NHCO}(\text{CH}_2)_9\text{CH}_3$		
343	$\text{NHCO}(\text{CH}_2)_{10}\text{CH}_3$		
344	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
345	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
346	$\text{NHCO}(\text{CH}_2)_{12}\text{CH}_3$		
347	$\text{NHCO}(\text{CH}_2)_7\text{CH}_3$		
348	$\text{NHCO}(\text{CH}_2)_7\text{CH}_3$		
349	$\text{NHCO}(\text{CH}_2)_7\text{CH}_3$		
350			
351	$\text{NHCO}(\text{CH}_2)_{11}\text{CH}_3$		
352	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		
355	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		
356	$\text{NHCONH}(\text{CH}_2)_{10}\text{CH}_3$		

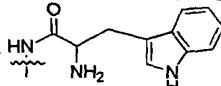
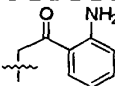
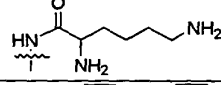
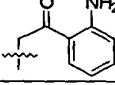
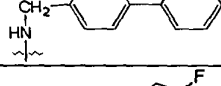
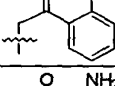
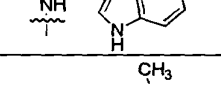
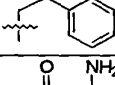
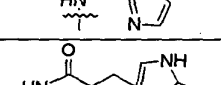
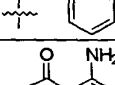
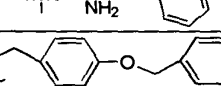
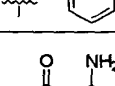
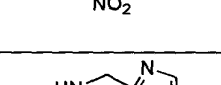
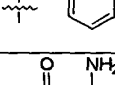
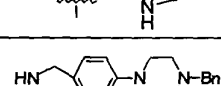
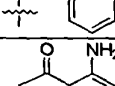
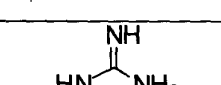
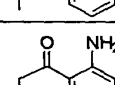
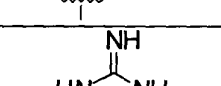
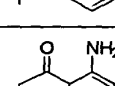
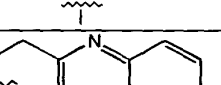
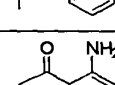
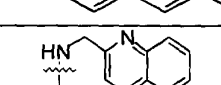
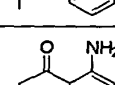
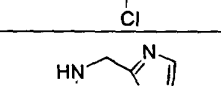
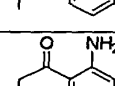
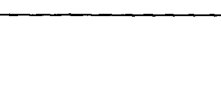
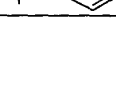
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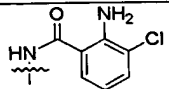
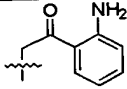
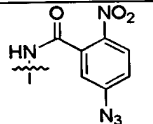
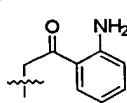
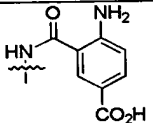
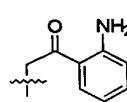
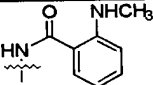
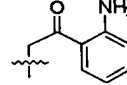
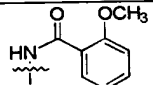
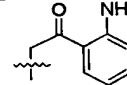
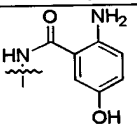
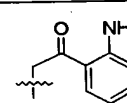
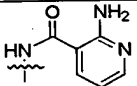
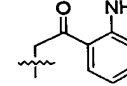
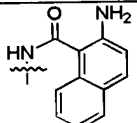
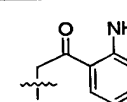
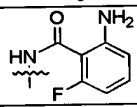
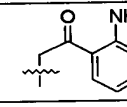
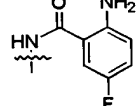
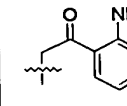
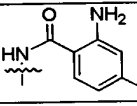
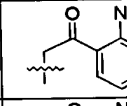
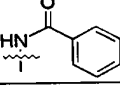
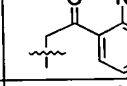
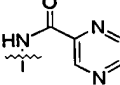
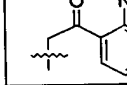
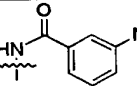
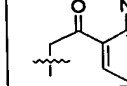
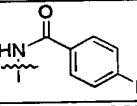
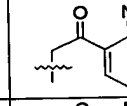
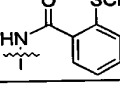
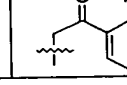
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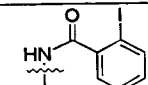
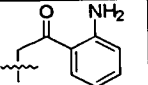
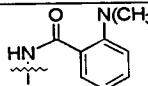
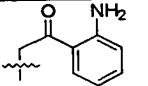
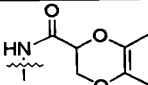
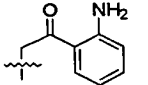
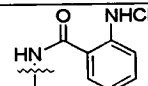
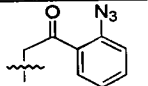
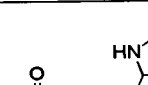
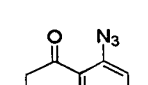
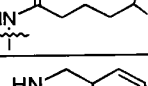
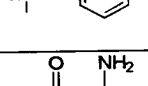
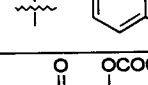
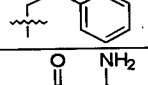
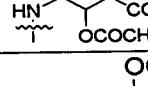
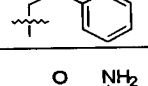
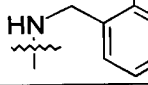
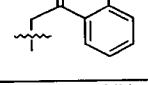
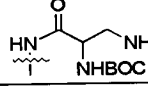
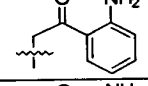
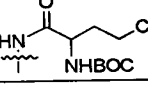
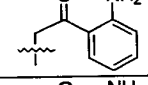
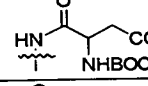
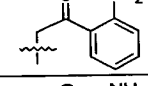
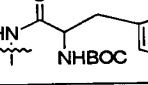
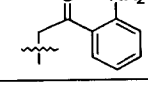
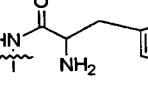
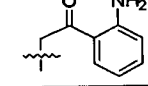
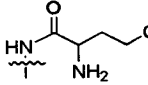
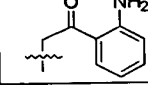
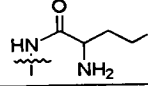
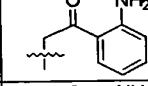
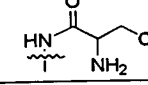
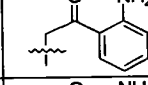
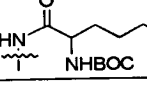
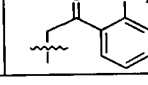
Cpd #	R	R ¹	R ²
45	NHCO(CH ₂) ₈ CH ₃		
54	NHCO(CH ₂) ₈ CH ₃		
76	NHCO(CH ₂) ₈ CH ₃		
81	NHCO(CH ₂) ₈ CH ₃		
85	NHCO(CH ₂) ₈ CH ₃		
102	NHCO(CH ₂) ₁₁ CH ₃		
209	NHCO(CH ₂) ₈ CH ₃		
212	NHCO(CH ₂) ₈ CH ₃		
253	NHCO(CH ₂) ₈ CH ₃		
260	NHCO(CH ₂) ₉ CH ₃		
262	NHCO(CH ₂) ₁₂ CH ₃		
282	NHCO(CH ₂) ₈ CH ₃		
285	NHCO(CH ₂) ₈ CH ₃		
319	NHCO(CH ₂) ₉ CH ₃		

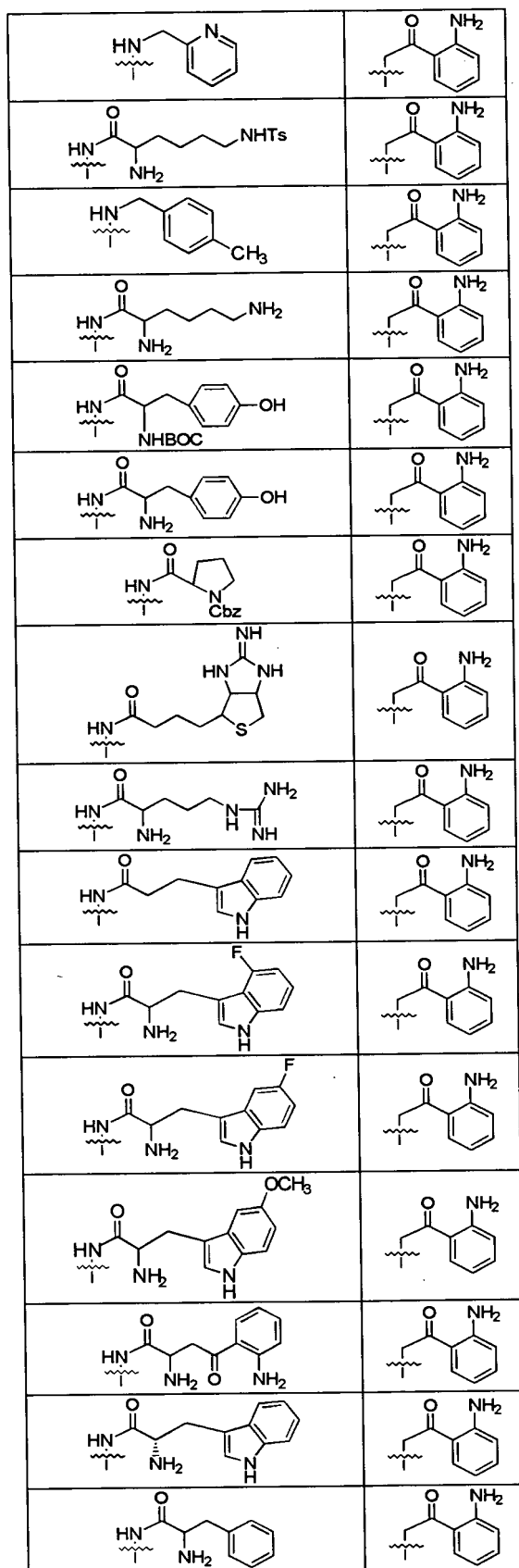
Cpd #	R	R ¹	R ²
322	NHCO(CH ₂) ₁₁ CH ₃		
333	NHCO(CH ₂) ₁₀ CH ₃		
334	NHCO(CH ₂) ₁₀ CH ₃		
335	NHCONH(CH ₂) ₁₁ CH ₃		
336	NHCONH(CH ₂) ₁₁ CH ₃		
344	NHCO(CH ₂) ₁₂ CH ₃		
355	NHCONH(CH ₂) ₁₀ CH ₃		

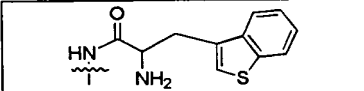
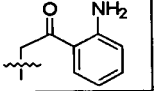
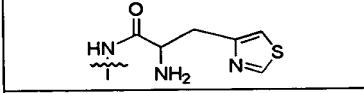
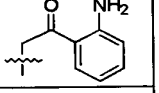
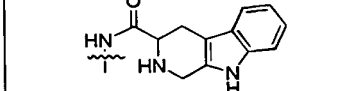
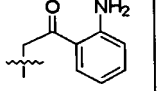
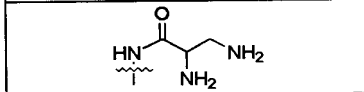
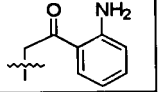
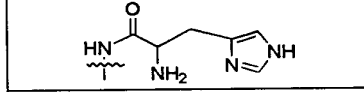
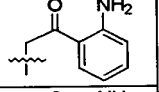
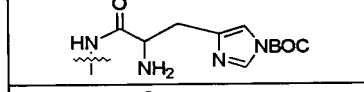
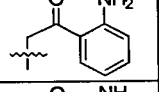
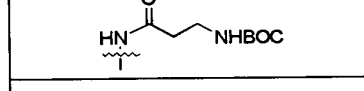
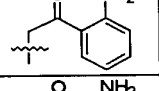
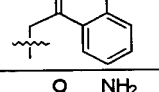
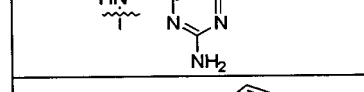
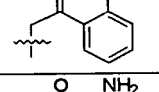
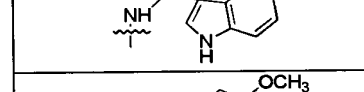
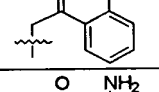
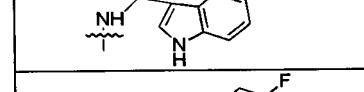
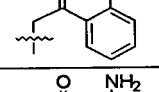
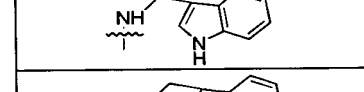
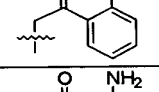
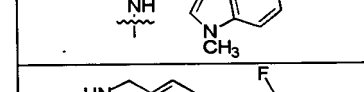
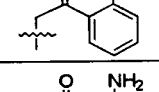
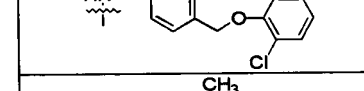
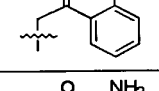
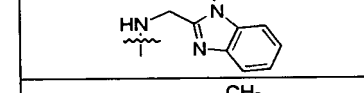
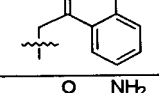
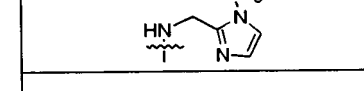
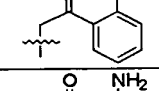
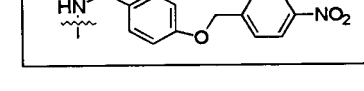
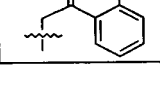
15. A compound of formula (I) according to any one of claims 1-4, wherein R is NHCO-[(C₆-C₁₄)-alkyl]-CH₃, and R¹ and R² are selected from:

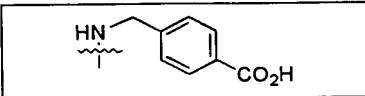
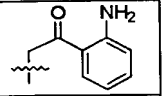
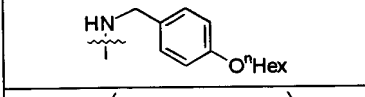
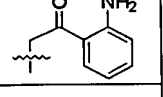
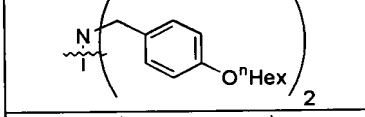
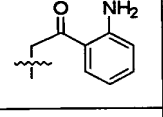
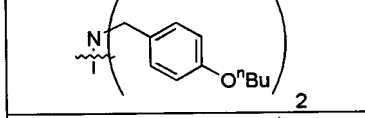
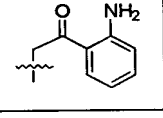
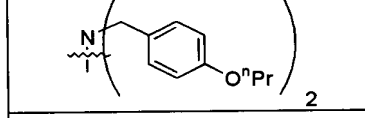
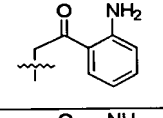
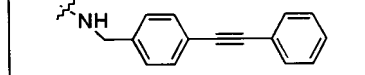
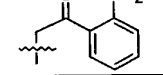
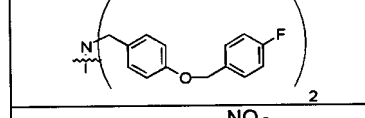
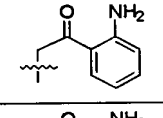
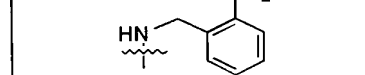
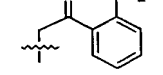
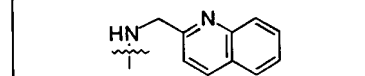
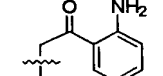
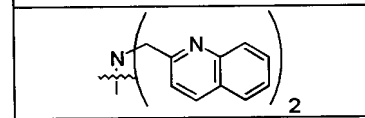
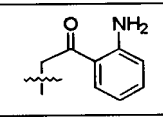
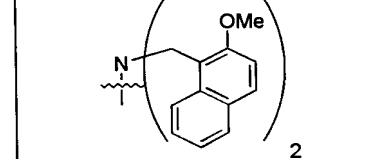
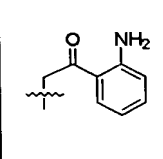
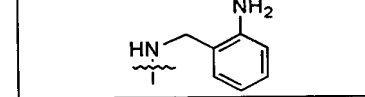
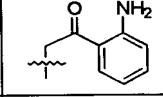
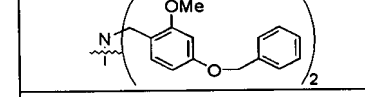
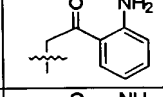
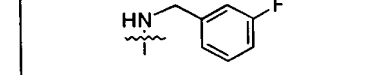
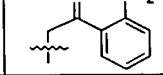
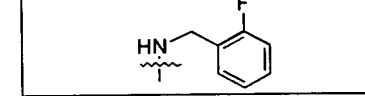
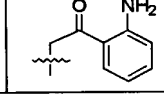
R ¹	R ²
NHSO ₂ Ph	

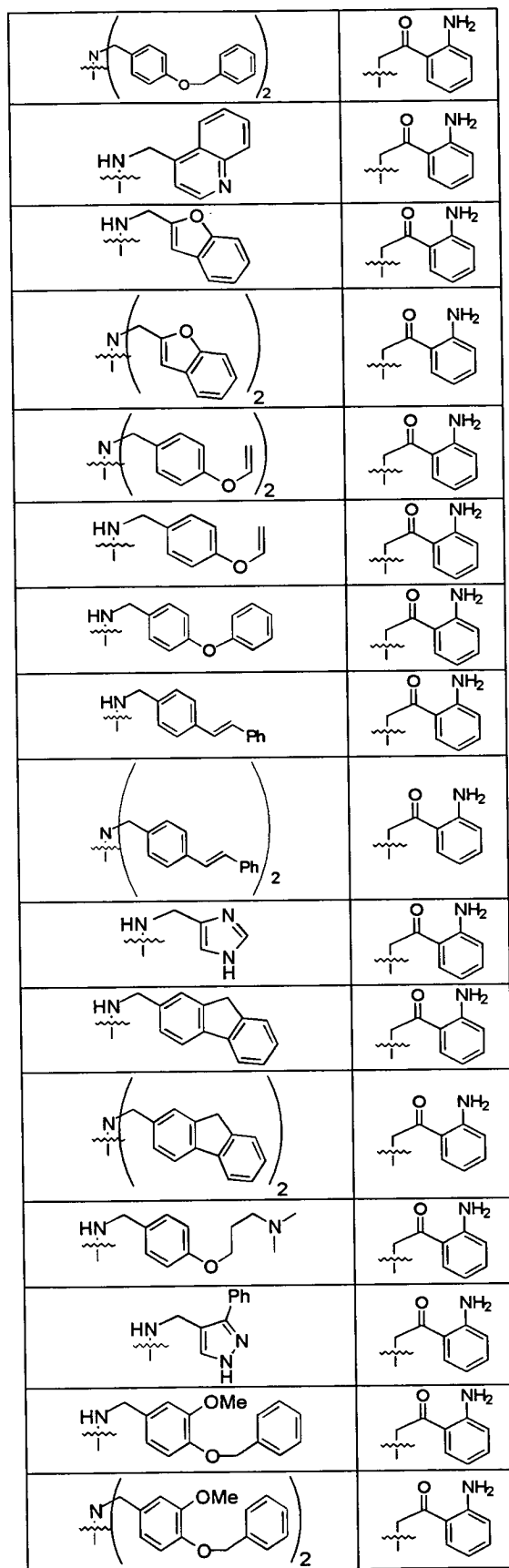
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

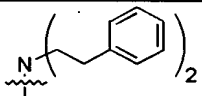
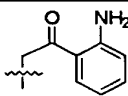
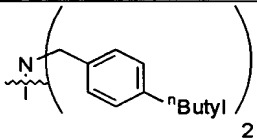
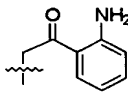
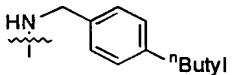
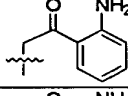
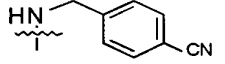
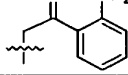
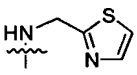
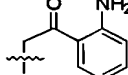
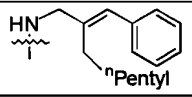
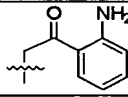
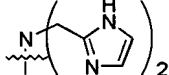
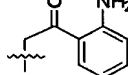
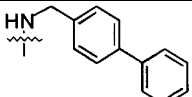
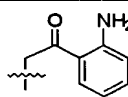
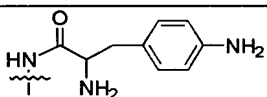
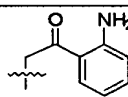
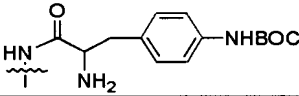
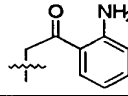
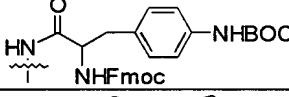
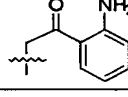
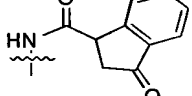
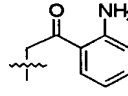
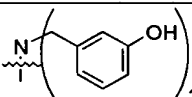
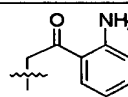
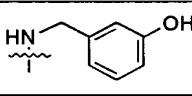
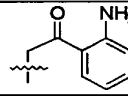
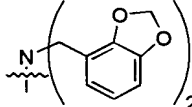
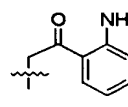
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

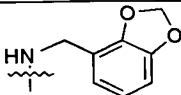
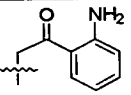
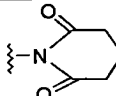
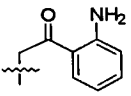
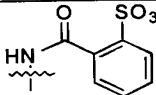
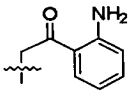
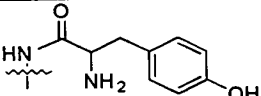
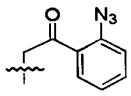
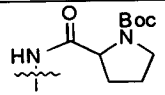
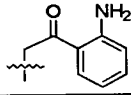
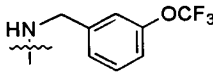
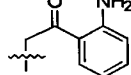
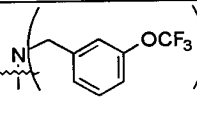
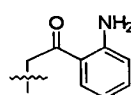
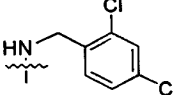
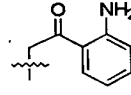
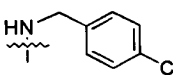
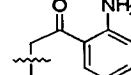
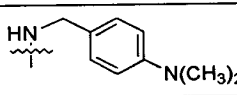
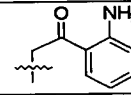
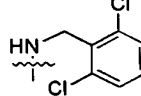
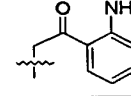
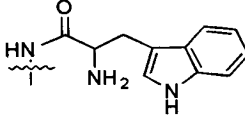
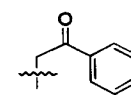
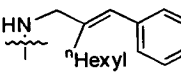
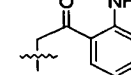
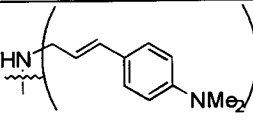
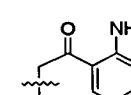
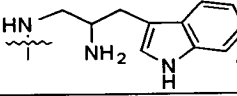
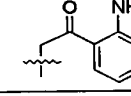
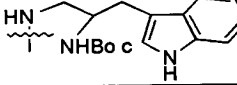
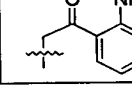


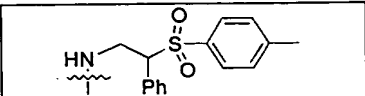
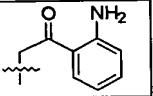
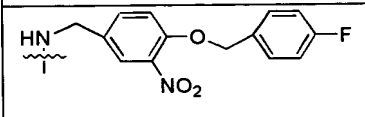
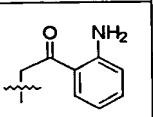
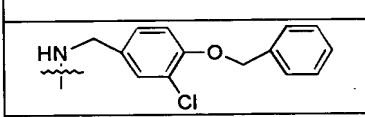
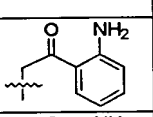
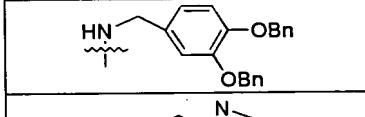
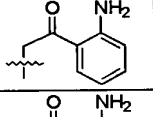
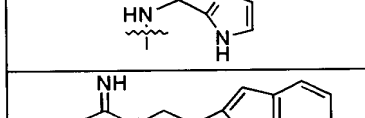
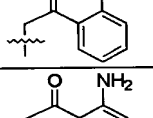
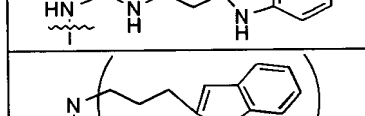
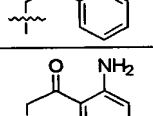
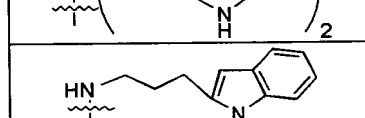
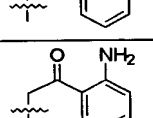
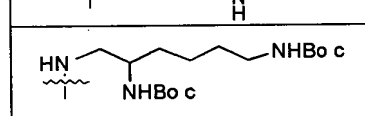
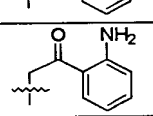
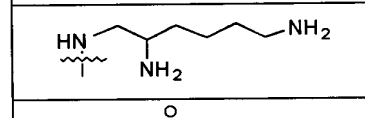
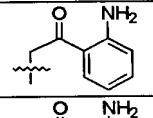
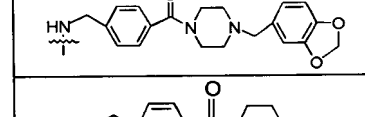
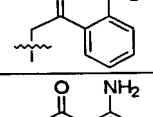
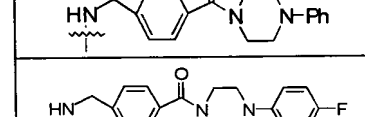
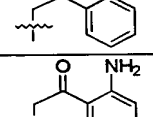
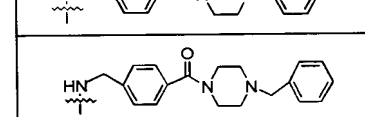
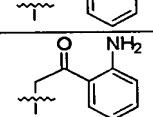
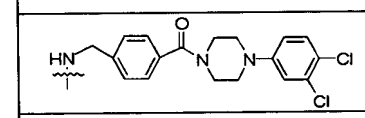
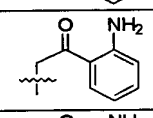
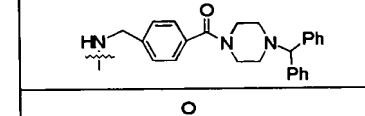
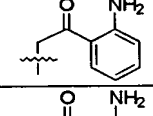
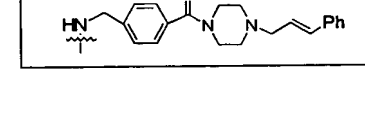
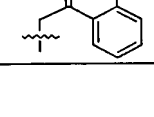




	
	
	
	
	
	
	
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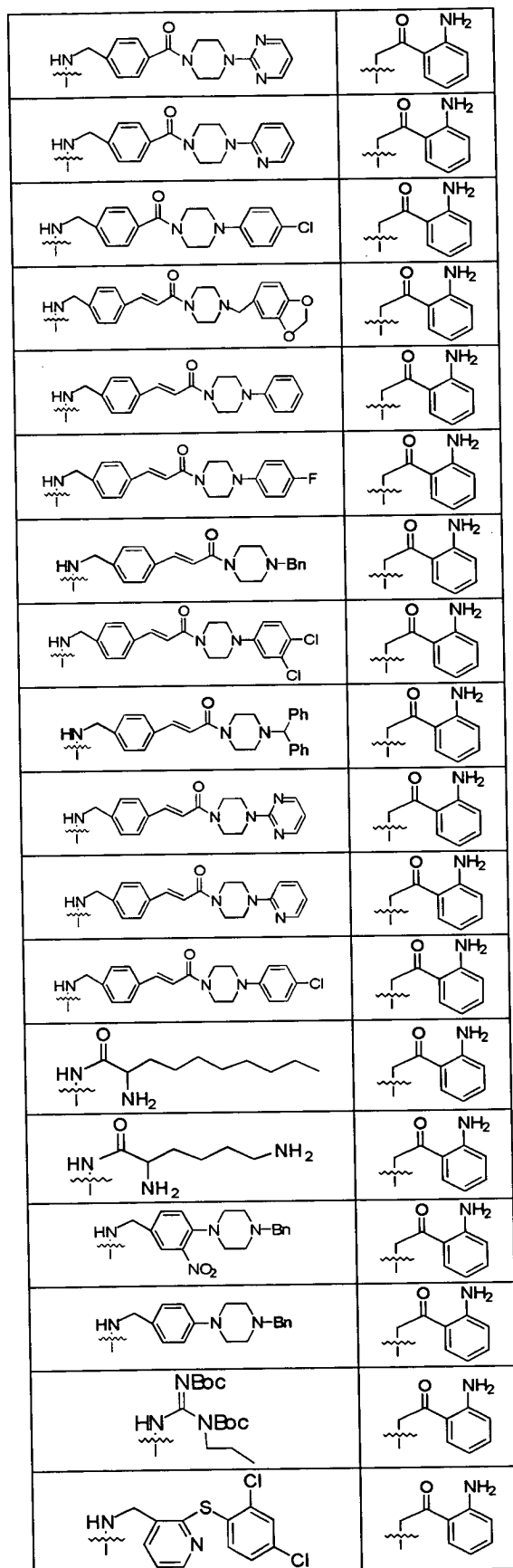
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

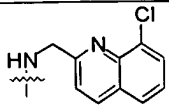
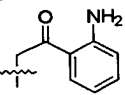
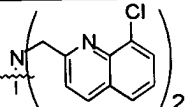
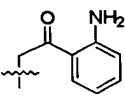
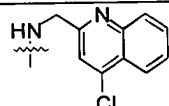
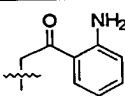
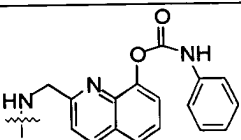
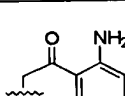
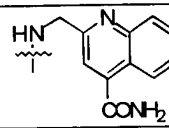
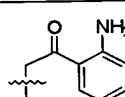
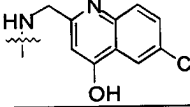
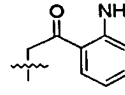
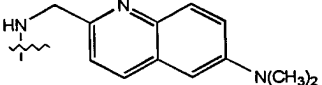
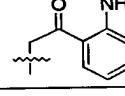
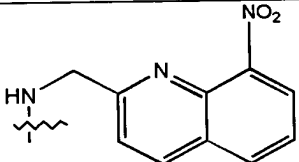
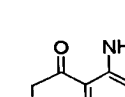
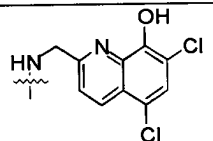
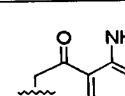
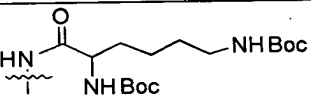
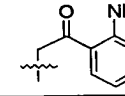
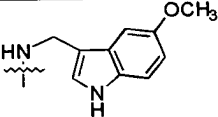
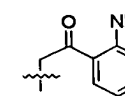
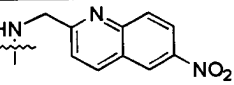
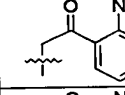
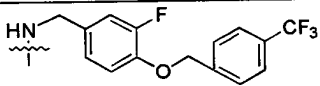
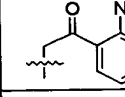
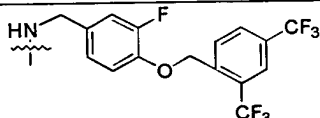
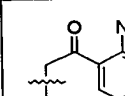


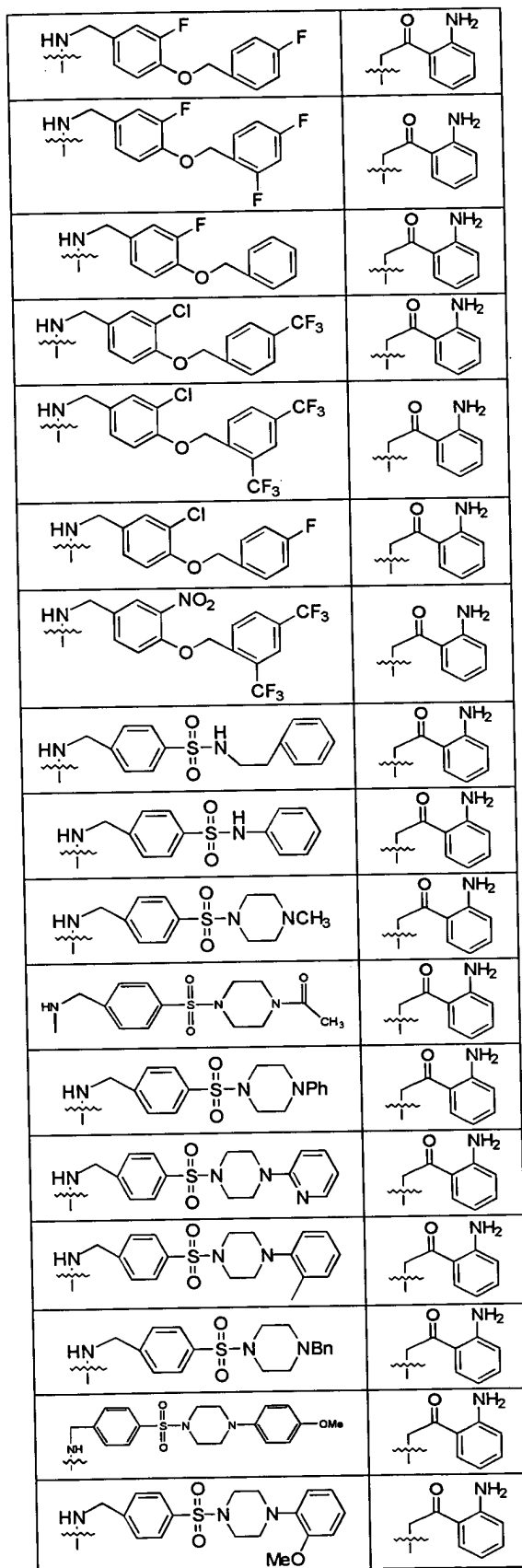
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

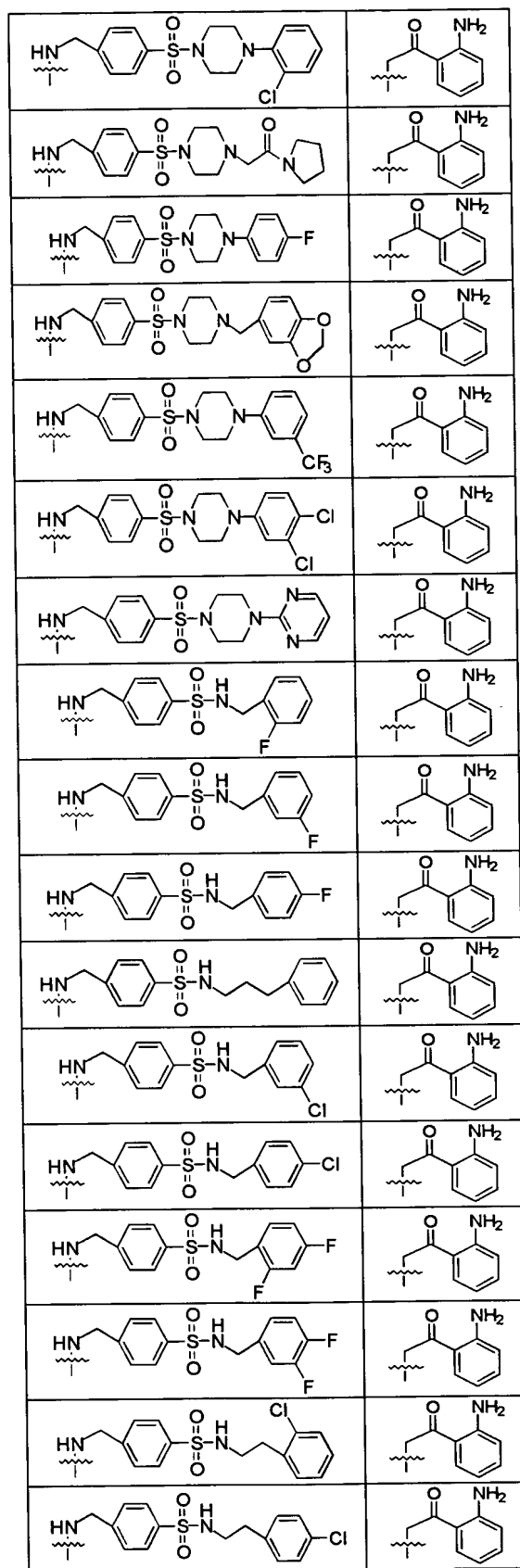
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

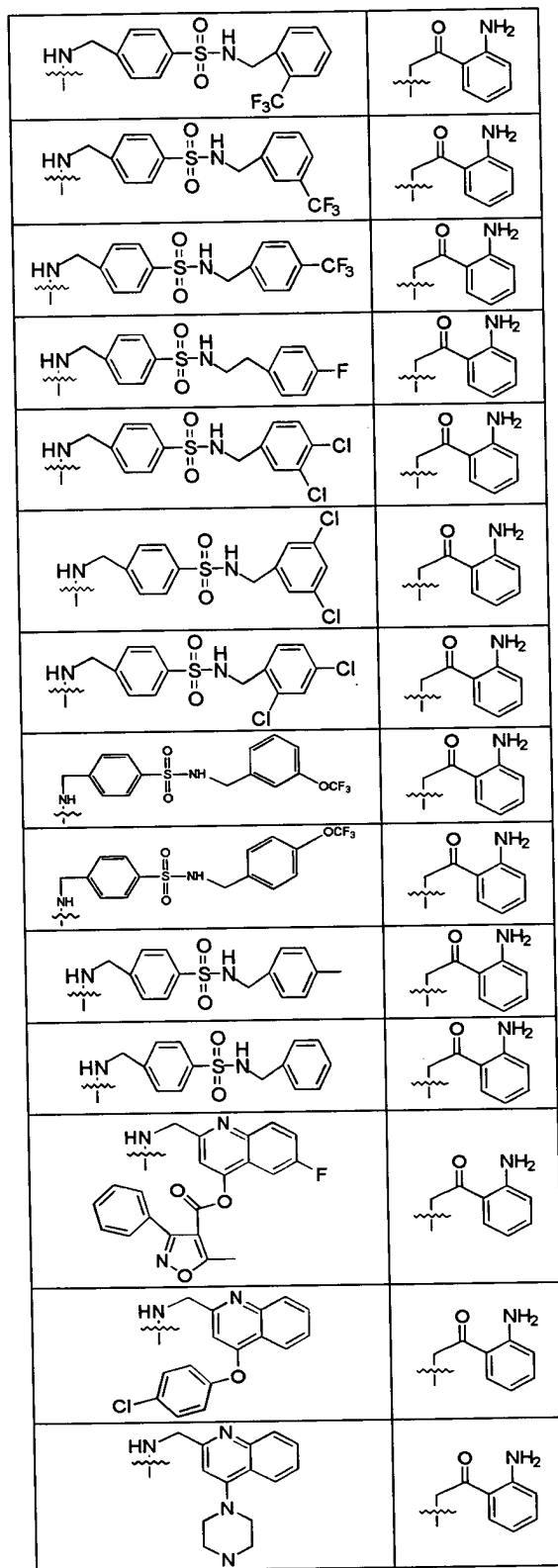
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	
	

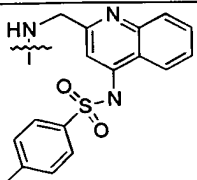
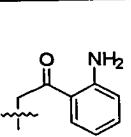
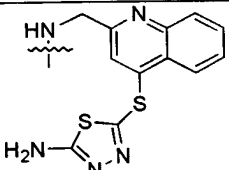
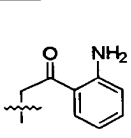
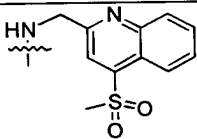
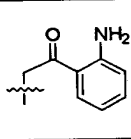
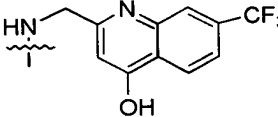
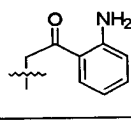
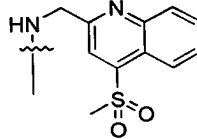
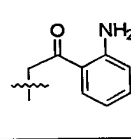
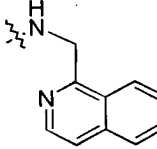
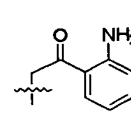
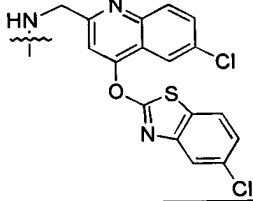
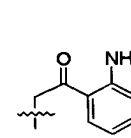
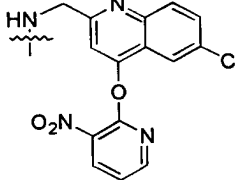
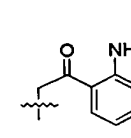
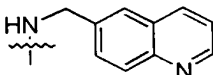
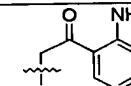
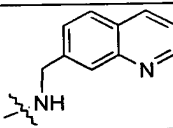
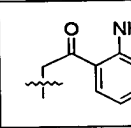








16. The compound according to claim 15, wherein R is selected from $\text{NHCO}[(\text{CH}_2)_{6-14}]\text{CH}_3$.
17. A pharmaceutical composition comprising the compound according to any one of claims 1-4 and a pharmaceutically acceptable carrier.
18. A method of treating a bacterial infection in a subject, comprising the step of administering a therapeutically-effective amount of the pharmaceutical composition according to claim 17 to a subject in need thereof.
19. The method according to claim 18, wherein said subject is selected from the group consisting of a human, an animal, a cell culture or a plant.
20. The method according to claim 18, wherein said bacterial infection is caused by a gram-positive bacteria.
21. The method according to claim 20, wherein said bacteria is an antibiotic-resistant bacteria.
22. The method according to claim 21, wherein said antibiotic-resistant bacteria are resistant to an antibiotic selected from the group consisting of vancomycin, methicillin, glycopeptide antibiotics, penicillin and daptomycin.
23. The method according to claim 18, further comprising the step of co-administering more than one compound of Formula (I) to a subject in need thereof.
24. The method according to claim 18, further comprising the step of co-administering an antimicrobial agent other than a compound of Formula (I) to a subject in need thereof.
25. The method according to claim 22, wherein said antimicrobial agent is selected from the group consisting of penicillins and related drugs,

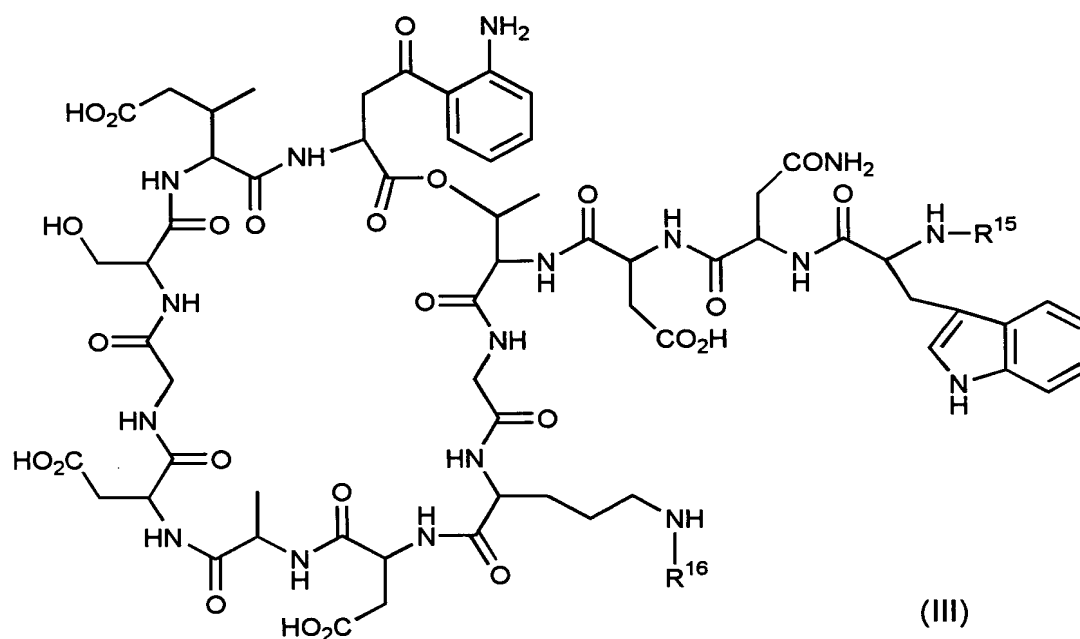
carbapenems, cephalosporins and related drugs, aminoglycosides, bacitracin, gramicidin, mupirocin, chloramphenicol, thiamphenicol, fusidate sodium, lincomycin, clindamycin, macrolides, novobiocin, polymyxins, rifamycins, spectinomycin, tetracyclines, vancomycin, teicoplanin, streptogramins, anti-folate agents including sulfonamides, trimethoprim and its combinations and pyrimethamine, synthetic antibacterials including nitrofurans, methenamine mandelate and methenamine hippurate, nitroimidazoles, quinolones, fluoroquinolones, isoniazid, ethambutol, pyrazinamide, para-aminosalicylic acid (PAS), cycloserine, capreomycin, ethionamide, prothionamide, thiacetazone, viomycin, eveminomycin, glycopeptide, glycylycylcline, ketolides, oxazolidinone; imipenen, amikacin, netilmicin, fosfomycin, gentamicin, ceftriaxone, Ziracin, LY 333328, CL 331002, HMR 3647, Linezolid, Synercid, Aztreonam, and Metronidazole, Epiroprim, OCA-983, GV-143253, Sanfetrinem sodium, CS-834, Biapenem, A-99058.1, A-165600, A-179796, KA 159, Dynemicin A, DX8739, DU 6681; Cefluprenam, ER 35786, Cefoselis, Sanfetrinem celexetil, HGP-31, Cefpirome, HMR-3647, RU-59863, Mersacidin, KP 736, Rifalazil; Kosan, AM 1732, MEN 10700, Lenapenem, BO 2502A, NE-1530, PR 39, K130, OPC 20000, OPC 2045, Venepirim, PD 138312, PD 140248, CP 111905, Sulopenem, ritipenam acoxyl, RO-65-5788, Cyclothialidine, Sch-40832, SEP-132613, micacocidin A, SB-275833, SR-15402, SUN A0026, TOC 39, carumonam, Cefozopran, Cefetamet pivoxil, and T 3811.

26. The method according to claim 22, wherein said antimicrobial agent is selected from the group consisting of imipenen, amikacin, netilmicin, fosfomycin, gentamicin, ceftriaxone, teicoplanin, Ziracin, LY333328, CL331022, HMR3647, Linezolid, Synercid, Aztreonam and Metronidazole.

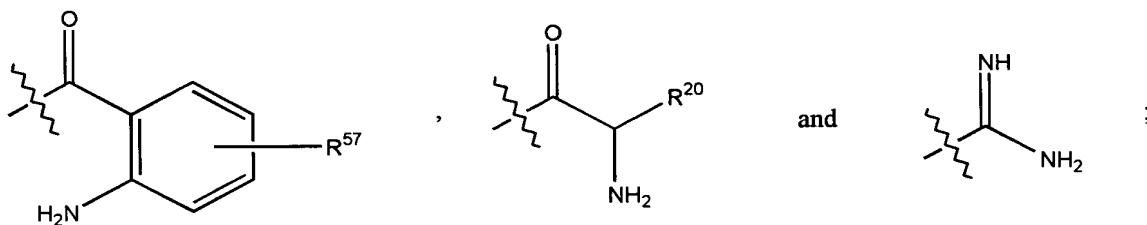
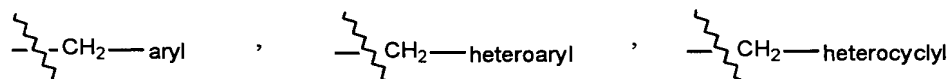
27. The method according to claim 19, wherein said subject is selected a human or an animal.

28. The method according to claim 27, wherein said subject is a human.

29. A compound having the formula (III):

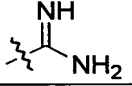
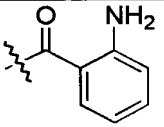
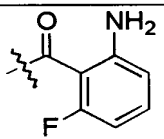
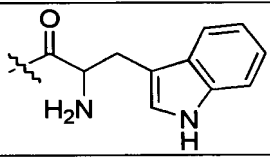
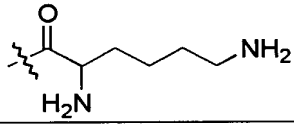
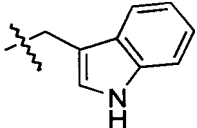
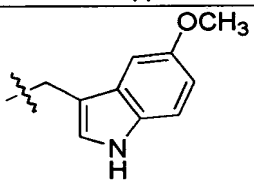
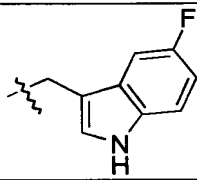
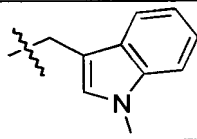
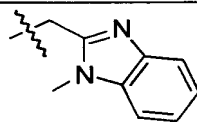


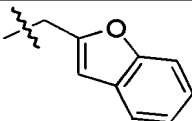
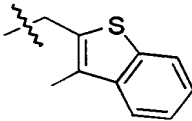
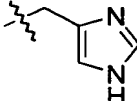
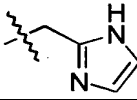
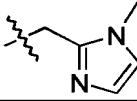
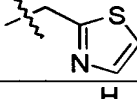
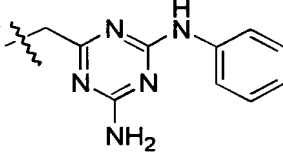
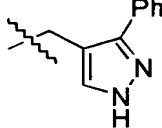
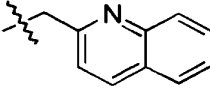
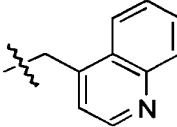
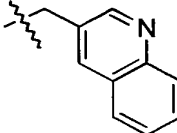
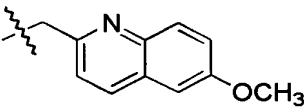
wherein R^{15} is selected from hydrido and an carbamate amino protecting group, preferably a *tert*-butoxycarbonyl group; wherein R^{16} is selected from the group consisting of

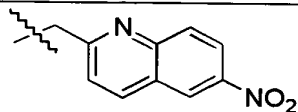
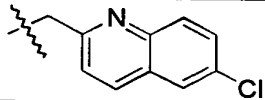
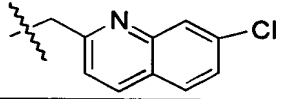
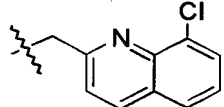
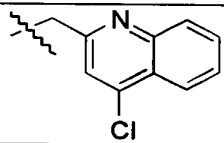
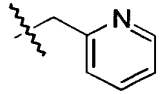
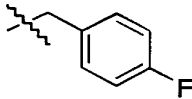
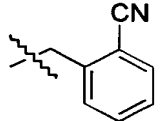
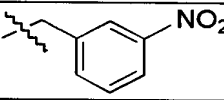
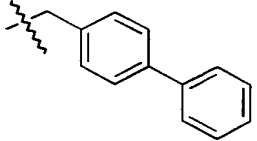
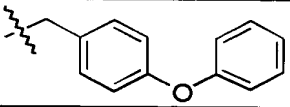
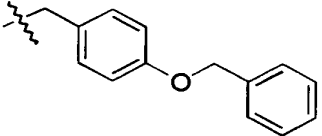


wherein R^{57} is a halo or halo substituted alkyl group, preferably a fluoro or trifluoromethyl group; wherein, R^{20} is an amino acid side chain, preferably a lysine or tryptophan side chain.

30. The compound according to claim 29 selected from:

Compound #	R ¹⁶
2	
10	
25	
45	
54	
79	
80	
81	
82	
84	

Compound #	R ¹⁶
139	
158	
146	
212	
85	
174	
78	
150	
130	
138	
168	
274	

Compound #	R ¹⁶
317	
280	
275	
283	
285	
50	
38	
115	
105	
76	
147	
164	

Compound #	R^{16}
210	
107	
111	
103	
253	
227	
372	
386	

31. A method of using the compound according to either of claims 29 or 30 to make a compound according to any one of claims 1-4.